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REGION III  
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**SUBJECT:** Risk-Based Concentration Table

**DATE:** 5/8/2001

**FROM:** Jennifer Hubbard, Toxicologist  
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**TO:** RBC Table Users

Attached is the EPA Region III Risk-Based Concentration (RBC) Table, which we prepare and post periodically for all interested parties.

**IMPORTANT NOTES:** To make the RBC Table more accessible and to minimize paper usage, it is now primarily available through the Internet. The address is <http://www.epa.gov/reg3hwmd/risk/riskmenu.htm>. The Table is available in both Lotus and Excel as "self-extracting" files. These files should be downloaded and then processed with your computer's "run" function. The files can then be viewed in Lotus or Excel. If you have technical questions about the toxicological or risk assessment aspects of the RBCs, please contact Jennifer Hubbard at 215-814-3328 or [hubbard.jennifer@epa.gov](mailto:hubbard.jennifer@epa.gov). Other questions can be addressed to Terri Fields at 215-814-3041. You can also consult the Frequently Asked Questions, below.

**SPECIAL NOTES ON THE MAY 2001 VERSION OF THE RBC TABLE**

The Region III RBC Table is typically updated in April and October. Users of the Table may have noticed the website posting explaining that an April release was not feasible this year. As the posting indicated, changes to chlorine dioxide, 1,3-dichlorobenzene, and vinyl chloride were anticipated. Cobalt has also been changed, due to the recent receipt of a new provisional RfD and RfC.

There are substantial changes to the vinyl chloride methodology. Readers are referred to the memo, "Derivation of Vinyl Chloride RBCs," which is a companion document to this RBC Table.

**CONTENTS, USES, AND LIMITATIONS OF THE RBC TABLE**

The RBC Table contains Reference Doses (RfDs) and Cancer Slope Factors (CSFs) for 400-500 chemicals. These toxicity factors have been combined with "standard" exposure scenarios to calculate RBCs--chemical concentrations corresponding to fixed levels of risk (i.e., a

Hazard Quotient (HQ) of 1, or lifetime cancer risk of 1E-6, whichever occurs at a lower concentration) in water, air, fish tissue, and soil.

The Region III toxicologists use RBCs to screen sites not yet on the NPL, respond rapidly to citizen inquiries, and spot-check formal baseline risk assessments. The primary use of RBCs is for chemical screening during baseline risk assessment (see EPA Regional Guidance EPA/903/R-93-001, "Selecting Exposure Routes and Contaminants of Concern by Risk-Based Screening"). The exposure equations come from EPA's Risk Assessment Guidance for Superfund (RAGS), while the exposure factors are those recommended in RAGS or supplemental guidance from the Superfund program. The attached technical background document provides specific equations and assumptions. Simply put, RBCs are like risk assessments run in reverse. For a single contaminant in a single medium, under standard default exposure assumptions, the RBC corresponds to the target risk or hazard quotient.

RBCs also have several important limitations. Specifically excluded from consideration are (1) transfers from soil to air, (2) cumulative risk from multiple contaminants or media, and (3) dermal risk. Additionally, the risks for inhalation of vapors from water are based on a very simple model, whereas detailed risk assessments may use more detailed showering models. Also, the toxicity information in the Table has been assembled by hand and (despite extensive checking and years of use) may contain errors. It's advisable to cross-check before relying on any RfDs or CSFs in the Table. If you note any errors, please let us know.

It is important to note that this Table uses inhalation RfDs and CSFs rather than RfCs (Reference Concentrations) and inhalation unit cancer risks. This is because the latter factors incorporate exposure assumptions and therefore can only be used for one exposure scenario. Because risk assessors need to evaluate risks for many types of scenarios, the factors have been converted to the more traditional RfDs and CSFs. Unless otherwise indicated in the toxicity-factor source, the assumption is that RfCs and unit risks should be adjusted by a 70-kilogram body weight and a 20 m<sup>3</sup>/day inhalation rate to generate the RfDs and CSFs.

Many users want to know if the RBCs can be used as valid no-action levels or cleanup levels, especially for soils. The answer is a bit complex. First, it is important to realize that the RBC Table does not constitute regulation or guidance, and should not be viewed as a substitute for a site-specific risk assessment. For sites where:

1. A single medium is contaminated;
2. A single contaminant contributes nearly all the health risk;
3. Volatilization, dermal contact, and other pathways not included in the RBCs are not expected to be significant;
4. The exposure scenarios and assumptions used in the RBC table are appropriate for the site;

- 5. The fixed risk levels used in the RBC table are appropriate for the site; and
- 6. Risk to ecological receptors is not expected to be significant;

the RBCs would probably be protective as no-action levels or cleanup goals. However, to the extent that a site deviates from this description, as most do, the RBCs would not necessarily be appropriate.

To summarize, the Table should generally not be used to set cleanup or no-action levels at CERCLA sites or RCRA Corrective Action sites, to substitute for EPA guidance for preparing baseline risk assessments, or to determine if a waste is hazardous under RCRA.

#### SPECIAL NOTES

The RBC Table was originally developed by Roy L. Smith, Ph.D., for use by risk assessors in the Region III Superfund program. Dr. Smith is no longer with Region III, and the Table continues to evolve. You may notice some modifications of formatting and conventions used in the Table.

For instance, besides formatting, the following changes are noteworthy:

As usual, updated toxicity factors have been used wherever available. However, because IRIS and provisional values are updated more frequently than the RBC Table, RBC Table users are ultimately responsible for obtaining the most up-to-date values. The RBC Table is provided as a convenience, but toxicity factors are compiled from the original sources and it is those original sources that should serve as the definitive reference.

Certain outdated and withdrawn numbers have been removed from the Table.

Changes to the table since the last semi-annual version have been marked with asterisks (\*\*). Changes may involve a corrected CAS number or a correction in the VOC status, a change in the SSL, or changes of RfDs and CSFs on IRIS.

RBCs are no longer rounded to 1E6 ppm. For certain low-toxicity chemicals, the RBCs exceed possible concentrations at the target risks. In such cases, Dr. Smith rounded these numbers to the highest possible concentration, or 1E6 ppm. This type of truncation has been discontinued so that Table users can adjust the RBCs to a different target risk whenever necessary. For example, when screening chemicals at a target HQ of 0.1, noncarcinogenic RBCs may simply be divided by 10. Such scaling is not possible when RBCs are rounded. Users who are interested in truncation can also consult the Soil Screening Guidance for a discussion of "C<sub>sat</sub>," the saturation concentration.

This Table was originally compiled to assist Superfund risk assessors in screening hazardous waste sites. The large number of chemicals made the Table unwieldy and

difficult to keep current. Many of the chemicals did not typically (or even occasionally) appear at Superfund sites. Starting with the April 1998 version of the Table, the 600+ chemicals were reduced to some 400-500 chemicals by eliminating many of those atypical chemicals. Through time, the Table may continue to grow or decrease in size. Comments on this issue are appreciated. During the last two and a half years, only one request was received for restoration of a chemical: NuStar was restored to the Table. (A list of the deleted chemicals is attached.)

At Region III Superfund sites, noncancer RBCs are typically adjusted downward to correspond to a target HQ of 0.1 rather than 1. (This is done to ensure that chemicals with additive effects are not prematurely eliminated during screening. Note that the RBCs displayed on the table are shown at an HQ of 1; to arrive at the RBC at 0.1, data users must do the conversion themselves.) However, some chemicals have RBCs at HQs of 0.1 that are lower than their RBCs at 1E-6 cancer risk. In other words, the screening RBC would change from carcinogenic to noncarcinogenic. A feature of this Table is that these chemicals are now flagged with a “!” symbol. Therefore, assessors screening with adjusted RBCs will be alerted to this situation.

Earlier versions of this Table included a substitution of inhalation toxicity factors for oral factors whenever oral factors were unavailable (this applied only to groundwater and air, but not soil or fish). This practice has been discontinued in order to minimize the uncertainty associated with such a conversion. The discontinuation of this practice does not significantly decrease the number of available RBCs.

The criterion for “VOC status” is in accordance with RAGS Part B: chemicals with Henry’s Law constants greater than 1E-5 and molecular weight less than 200 are now marked as VOCs. This increases consistency with the national guidance and with other EPA regions that use risk-based screening numbers.

Earlier versions of this Table included soil screening levels (SSLs), when those values were available in draft form. Since the finalization of the SSL Guidance, risk assessors are urged to consult the final SSL Guidance directly. However, for generic use in Region III, the table now contains soil-to-groundwater SSLs in accordance with the new guidance. For more information, see the Region III memo on SSLs, or consult the national SSL guidance directly (Soil Screening Guidance: User’s Guide, April 1996, Publication 9355.4-23; and Soil Screening Guidance: Technical Background Document, May 1996; EPA/540/R-95/128).

You may notice there are now two rows for uranium, one reflecting the IRIS (EPA consensus) value and the other reflecting a more recent, but provisional value. Region III has shown both on this table, rather than choosing one over the other, to give Table users as much information as possible.

## FREQUENTLY ASKED QUESTIONS

To help you better understand the RBC Table, here are answers to our most often-asked questions:

1. How can the age-adjusted inhalation factor (11.66) be less than the inhalation rate for either a child (12) or an adult (20)?

Age-adjusted factors are not intake rates, but rather partial calculations which have different units from intake rates. (Therefore, they are not directly comparable.) The fact that these partial calculations have values similar to intake rates is really coincidental, an artifact of the similar magnitude of years of exposure and time-averaged body weight.

2. For manganese, IRIS shows an oral RfD of 0.14 mg/kg/day, but the RBC Table uses 2E-2 mg/kg/day. Why?

The IRIS RfD includes manganese from all sources, including diet. The explanatory text in IRIS recommends using a modifying factor of 3 when calculating risks associated with non-food sources, and the Table follows this recommendation. IRIS also recommends subtracting dietary exposure (default assumption in this case 5 mg). Thus, the IRIS RfD has been lowered by a factor of 2 x 3, or 6. The Table now reflects manganese RBCs for both “food” and “non-food” (most environmental) sources.

3. What is the source of the child’s inhalation rate of 12 m<sup>3</sup>/day?

The calculation comes from basic physiology. It’s a scaling of the mass-specific 20 m<sup>3</sup>/day rate for adults from a body mass of 70 kg to 15 kg, using the 2/3 power of mass, as follows:

$$\begin{aligned} \text{Ircm} &= \text{mass-specific child inhalation rate (m}^3/\text{kg/day)} \\ \text{Irc} &= \text{child inhalation rate (m}^3/\text{day)} \end{aligned}$$

$$20 \text{ m}^3/\text{day} / 70 \text{ kg} = 0.286 \text{ m}^3/\text{kg/day} \text{ (mass-specific adult inhalation rate)}$$

$$0.286 \text{ m}^3/\text{kg/day} \times (70^{0.67}) = (\text{Ircm}) \times (15^{0.67})$$

$$\text{Ircm} = 0.803 \text{ m}^3/\text{kg/day}$$

$$\text{Irc} = \text{Ircm} \times 15 \text{ kg} = 0.803 \text{ m}^3/\text{kg/day} \times 15 \text{ kg} = 12.04 \text{ m}^3/\text{day}$$

4. Can the oral RfDs in the RBC Table be applied to dermal exposure?

Not directly. Oral RfDs are usually based on administered dose and therefore tacitly

include a GI absorption factor. Thus, any use of oral RfDs (or CSFs) in dermal risk calculations should involve removing this absorption factor. Consult the Risk Assessment Guidance for Superfund, Part A, Appendix A, for further details on how to do this.

5. The exposure variables table in the RBC background document lists the averaging time for non-carcinogens as “ED\*365.” What does that mean?

ED is exposure duration, in years, and \* is the computer-ese symbol for multiplication. Multiplying ED by 365 simply converts the duration to days. In fact, the ED term is included in both the numerator and denominator of the RBC algorithms for non-cancer risk, canceling it altogether. See RAGS for more information.

6. Why is inorganic lead not included in the RBC Table?

EPA has no consensus RfD or CSF for inorganic lead, so it is not possible to calculate RBCs as we have done for other chemicals. EPA considers lead to be a special case because of the difficulty in identifying the classic “threshold” needed to develop an RfD.

EPA therefore evaluates lead exposure by using blood-lead modeling, such as the Integrated Exposure-Uptake Biokinetic Model (IEUBK). The EPA Office of Solid Waste has also released a detailed directive on risk assessment and cleanup of residential soil lead. The directive recommends that soil lead levels less than 400 mg/kg are generally safe for residential use. Above that level, the document suggests collecting data and modeling blood-lead levels with the IEUBK model. For the purposes of screening, therefore, 400 mg/kg is recommended for residential soils. For water, we suggest 15 ug/l (the EPA Action Level in water), and for air, the National Ambient Air Quality Standard.

7. Where did the CSFs for carcinogenic PAHs come from?

The PAH CSFs are all calculated relative to benzo[a]pyrene, which has an IRIS slope factor. The relative factors for the other PAHs can be found in “Provisional Guidance for Quantitative Risk Assessment of Polycyclic Aromatic Hydrocarbons,” Final Draft, ECAO-CIN-842 (March, 1993).

8. May I please have a copy of a previous RBC Table?

We do not distribute outdated copies of the RBC Table. Each new version of the Table supersedes all previous versions.

9. Please elaborate on the meaning of the “W” source code in the Table.

The “W” code means that a RfD or CSF is currently not present on either IRIS or HEAST, but that it was once present on either IRIS or HEAST and was removed. Such

withdrawal usually indicates that consensus on the number no longer exists among EPA scientists, but not that EPA believes the contaminant to be unimportant.

Withdrawn numbers are shown in the Table because we still need to deal with these contaminants during the long delays before replacement numbers are ready. For the purpose of screening, a "W" value is similar to a provisional value in that neither value has achieved Agency consensus. The "W" code should serve as a clear warning that before making any serious decision involving that contaminant, you will need to develop an interim value based on current scientific understanding.

If you are assessing risks at a site where a major contaminant is coded "W," consider working with your Region EPA risk assessor to develop a current toxicity constant. If the site is being studied under CERCLA, the EPA-NCEA Regional Technical Support group may be able to assist.

10. Can I get copies of supporting documents for interim toxicity constants which are coded "E" in the RBC Table?

Unfortunately, Region 3 does not have a complete set of supporting documents. The EPA-NCEA Superfund Technical Support Center prepares these interim toxicity constants in response to site-specific requests from Regional risk assessors, and sends the documentation only to the requestor. The RBC Tables contain only the latest interim values that we've either requested or have otherwise received. NCEA maintains the master data base of these chemicals, but will not release documentation of provisional values unless they are recent. Furthermore, since NCEA's Superfund Technical Support Center is mainly for the support of Superfund, it usually cannot develop new criteria unless authorized to do so for a specific Superfund project.

If an "E"-coded contaminant is a chemical of potential concern at your site, we urge you to work with the EPA Regional risk assessor assigned to the project in order to develop or obtain documentation for provisional values. EPA Region 3 furnishes documents only when needed to support Regional risk assessments or recommendations.

11. Why is there no oral RfD for mercury? How should I handle mercury?

IRIS gives oral RfDs for mercuric chloride and for methylmercury, but not for elemental mercury. Therefore, the RBC Table reflects this primary source. Consult your toxicologist to determine which of the available mercury numbers is suitable for the conditions at your site (e.g., whether mercury is likely to be organic or inorganic.)

Attachment

## “DISCONTINUED” CHEMICALS

These chemicals may still have toxicity criteria available in IRIS, HEAST, or NCEA provisional values. However, they are not routine chemicals and therefore will not be routinely maintained in the RBC Table, unless our Table users report a significant need for chemicals to be re-added. Some of the chemicals on this Table were deleted because supporting toxicity information has been withdrawn or is unavailable.

acephate	acetone cyanohydrin
aci fluorfen	acrylic acid
ally	allyl alcohol
aluminum phosphide	amdro
ametryn	m-aminophenol
amitraz	ammonium sulfamate
antimony potassium tartrate	apollo
aramite	asulam
avermectin B1	barium cyanide
bayleton	benefin
benomyl	benzotrichloride
bidrin	biphenthin
bis(2-chloro-1-methylethyl)ether	boron trifluoride
bisphenol A	bromoxynil
4-bromophenyl phenyl ether	butylphthalyl butylglycolate
bromoxynil octanoate	captanol
cacodylic acid	carboxin
captan	chlorimuron-ethyl
chloramben	2-chloroacetophenone
chloroacetaldehyde	2-chloroethyl vinyl ether
4-chlorobenzotrifluoride	
4-chloro-2-methylaniline hydrochloride	chlorpropham
chlorothalonil	chlorthiophos
chlorsulfuron	
coal tar creosote	
cyclohexlamine	cyromazine
danitol	decabromodiphenyl ether
demeton	diallate
diethylforamide	diflubenzuron
dimethipin	dimethoate
N,N-dimethylformamide	dimethyl terephthalate
diphenamid	direct black 38
direct blue 6	direct brown 95
dodine	1,2-epoxybutane
ethephon	2-ethoxyethanol acetate
ethyl acrylate	EPTC

ethylene cyanohydrin	
ethyl p-nitrophenyl phenylphosphorothioate	
ethylphthalyl ethyl glycolate	express
fluoridone	flurprimidol
flutolanil	fluvalinate
folpet	fostetyl-al
furium	furmecyclox
glufosinate-ammonium	haloxyfop-methyl
harmony	imazalil
imazaquin	iprodione
isoxaben	kepone
lactofen	linuron
londax	
maleic hydrazide	malononitrile
mancozeb	maneb
merphos	merphos oxide
metalaxy1	methamidophos
methomyl	2-methoxyethanol acetate
2-methoxyethanol	2-methoxy-5-nitroaniline
2-methylaniline hydrochloride	methyl chlorocarbonate
4,4-methylene bisbenzeneamine	metribuzin
molinate	2-naphthylamine
napropamide	
nickel subsulfide	nitrapyrin
3-nitroaniline	4-nitroaniline
nitroguanidine	norflurazon
octabromodiphenyl ether	
octamethylpyrophosphoramido	paclobutrazol
pebulate	pendimethalin
pentabromo-6-chlorocyclohexane	
pentabromodiphenyl ether	phenmedipham
phenylmercuric acetate	phorate
phosmet	picloram
pirimiphos-methyl	prochloraz
profluralin	pronamide
propargyl alcohol	propazine
propham	propiconazole
propylene oxide	pydrin
quinalphos	savey
selenourea	sethoxydim
sodium fluoroacetate	sodium metavanadate
systhane	tebuthiuron
temephos	terbacil
terbufos	terbutryn

tetrachlorovinphos	tetraethylidithiopyrophosphate
thallium selenide	
2-(thiocyanomethylthio)-benzothiazole	
thiofanox	thiophanate-methyl
thiram	tralomethrin
trallate	triasulfuron
2,4,6-trichloroaniline hydrochloride	
tridiphane	triethylamine
trifluralin	vernam

**EPA REGION III RISK-BASED CONCENTRATION TABLE:  
TECHNICAL BACKGROUND INFORMATION**

originally developed by Roy L. Smith, Ph.D., Toxicologist  
revised 4/12/99 by Jennifer Hubbard, Toxicologist

**Development of Risk-Based Concentrations**

**General**

Separate carcinogenic and non-carcinogenic risk-based concentrations were calculated for each compound for each pathway. The concentration in the table is the lower of the two, rounded to two significant figures. The following terms and values were used in the calculations:

Exposure variables	Value	Symbol
<i>General:</i>		
Carcinogenic potency slope oral (risk per mg/kg/d):	*	CPSo
Carcinogenic potency slope inhaled (risk per mg/kg/d):	*	CPSi
Reference dose oral (mg/kg/d):	*	RfDo
Reference dose inhaled (mg/kg/d):	*	RfDi
Target cancer risk:	1e-06	TR
Target hazard quotient:	1	THQ
Body weight, adult (kg):	70	BWa
Body weight, age 1-6 (kg):	15	BWc
Averaging time carcinogens (d):	25550	ATc
Averaging time non-carcinogens (d):	ED*365	ATn
Inhalation, adult (m <sup>3</sup> /d):	20	IRAA
Inhalation, child (m <sup>3</sup> /d):	12	IRAc
Inhalation factor, age-adjusted (m <sup>3</sup> -y/kg-d):	11.66	IFAadj
Tap water ingestion, adult (L/d):	2	IRWa
Tap water ingestion, age 1-6 (L/d):	1	IRWc
Tap water ingestion factor, age-adjusted (L-y/kg-d):	1.09	IFWadj
Fish ingestion (g/d):	54	IRF
Soil ingestion, adult (mg/d):	100	IRSa
Soil ingestion, age 1-6 (mg/d):	200	IRSc
Soil ingestion factor, age adjusted (mg-y/kg-d):	114.29	IFSadj
<i>Residential:</i>		
Exposure frequency (d/y):	350	EFr

Exposure variables	Value	Symbol
Exposure duration, total (y):	30	EDtot
Exposure duration, age 1-6 (y):	6	EDc
Volatilization factor (L/m3):	0.5	K
<i>Occupational:</i>		
Exposure frequency (d/y):	250	EFo
Exposure duration (y):	25	EDo
Fraction of contaminated soil ingested (unitless)	0.5	FC

\*: Contaminant-specific toxicological constants. The priority among sources of toxicological constants was as follows: (1) IRIS, (2) HEAST, (3) HEAST alternative method, (4) EPA-NCEA Superfund Health Risk Technical Support Center, (5) withdrawn from IRIS or HEAST, and (6) other EPA documents. Each source was used only if numbers from higher-priority sources were unavailable, unless NCEA indicated a newer provisional value was superior to an older HEAST value. The EPA Superfund Health Risk Technical Support Center, part of the EPA National Center for Environmental Assessment in Cincinnati, develops provisional RfDs and CPSs on request for contaminants not in IRIS or HEAST. These provisional values are labeled "E = EPA-NCEA provisional" in the table. It is possible they may be obsolete. If one of the "E" constants is important to a Superfund risk assessment, consider requesting, through a Regional risk assessor, a new provisional value.

### Age-adjusted factors

Because contact rates with tap water, ambient air, and residential soil are different for children and adults, carcinogenic risks during the first 30 years of life were calculated using age-adjusted factors. These factors approximated the integrated exposure from birth until age 30 by combining contact rates, body weights, and exposure durations for two age groups - small children and adults. The age-adjusted factor for soil was obtained from RAGS IB; the others were developed by analogy.

#### (1) Air inhalation

$$IFA_{adj} \frac{m^3 \cdot y}{kg \cdot d} = \frac{EDC \cdot IRAc}{BWC} + \frac{(EDtot - EDC) \cdot IRAa}{BWa}$$

#### (2) Tap water ingestion

$$IFW_{adj} \frac{L \cdot y}{kg \cdot d} = \frac{EDC \cdot IRWC}{BWC} + \frac{(EDtot - EDC) \cdot IRWa}{BWa}$$

#### (3) Soil ingestion

$$IFS_{adj} \frac{mg \cdot y}{kg \cdot d} = \frac{EDC \cdot IRS_c}{BWC} + \frac{(EDtot - EDC) \cdot IRS_a}{BWa}$$

## Residential water

Volatilization terms were calculated only for compounds with a mark in the "VOC" column. Compounds having a Henry's Law constant greater than  $10^{-5}$  and a molecular weight less than 200 were considered volatile. The list may be incomplete, but is unlikely to include false positives. The equations and the volatilization factor (K, above) were obtained from RAGS IB. Oral potency slopes and reference doses were used for both oral and inhaled exposures for volatile compounds lacking inhalation values. RBCs for carcinogens were based on combined childhood and adult exposure; for non-carcinogens RBCs were based on adult exposure.

### (4) Carcinogens

$$RBC \frac{\mu g}{L} = \frac{TR \cdot ATc \cdot 1000 \frac{\mu g}{mg}}{EFr \cdot ([K \cdot IFAadj \cdot CPSi] + [IFWadj \cdot CPSo])}$$

### (5) Non-carcinogens

$$RBC \frac{\mu g}{L} = \frac{THQ \cdot BWa \cdot ATn \cdot 1000 \frac{\mu g}{mg}}{EFr \cdot EDtot \cdot \left( \frac{K \cdot IRAa}{RfDi} + \frac{IRWa}{RfDo} \right)}$$

## Ambient air

Oral potency slopes and references were used where inhalation values were not available. RBCs for carcinogens were based on combined childhood and adult exposure; for non-carcinogens RBCs were based on adult exposure.

### (6) Carcinogens

$$RBC \frac{\mu g}{m^3} = \frac{TR \cdot ATc \cdot 1000 \frac{\mu g}{mg}}{EFr \cdot IFAadj \cdot CPSi}$$

### (7) Non-carcinogens

$$RBC \frac{\mu g}{m^3} = \frac{THQ \cdot RfDi \cdot BWa \cdot ATn \cdot 1000 \frac{\mu g}{mg}}{EFr \cdot EDtot \cdot IRAa}$$

## Edible fish

All RBCs were based on adult exposure.

### (8) Carcinogens

$$RBC \frac{mg}{kg} = \frac{TR \cdot BWa \cdot ATc}{EFr \cdot EDtot \cdot \frac{IRF}{1000 \frac{g}{kg}} \cdot CPSo}$$

### (9) Non-carcinogens

$$RBC \frac{mg}{kg} = \frac{THQ \cdot RfDo \cdot BWa \cdot ATn}{EFr \cdot EDtot \cdot \frac{IRF}{1000 \frac{g}{kg}}}$$

## Commercial/industrial soil ingestion

RBCs were based on adult occupational exposure, including an assumption that only 50% of total soil ingestion is work-related.

### (10) Carcinogens

$$RBC \frac{mg}{kg} = \frac{TR \cdot BWa \cdot ATc}{EFO \cdot EDO \cdot \frac{IRSA}{10^6 \frac{mg}{kg}} \cdot FC \cdot CPSo}$$

### (11) Non-carcinogens

$$RBC \frac{mg}{kg} = \frac{THQ \cdot RfDo \cdot BWa \cdot ATn}{EFO \cdot EDO \cdot \frac{IRSA}{10^6 \frac{mg}{kg}} \cdot FC}$$

## Residential soil ingestion

RBCs for carcinogens were based on combined childhood and adult exposure; RBCs for non-carcinogens were based on childhood exposure only.

### (12) Carcinogens

$$RBC \frac{mg}{kg} = \frac{TR \cdot ATc}{EFr \cdot \frac{IFSAadj}{10^6 \frac{mg}{kg}} \cdot CPSo}$$

## (13) Non-carcinogens

$$RBC \frac{mg}{kg} = \frac{THQ \cdot RfDo \cdot BWc \cdot ATn}{EFr \cdot EDc \cdot \frac{IRSc}{10^6 \frac{mg}{kg}}}$$



Sources: I = IRIS H = HEAST A = HEAST Alternate W = Withdrawn from IRIS or HEAST  
 E = EPA-NCEA provisional value O = other

Chemical	CAS	RfDo	CSFo	RfDI	CSFI	VOC	Basis: C = Carcinogenic effects N = Noncarcinogenic effects I = RBC at HI<0.1 < RBC-C					Region III SSLs		
		mg/kg/d	1/mg/kg/d	mg/kg/d	1/mg/kg/d		Tap water ug/l	Ambient air ug/m3	Fish mg/kg	Soil Industrial mg/kg	Residential mg/kg	DAF 1 mg/kg	DAF 20 mg/kg	
ACETALDEHYDE	75070			2.57E-003 I	7.7E-003 I	y	1.6E+000 C	8.1E-001 C						
ACETOCHLOR	34256821	2E-002 I					7.3E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N	3.8E-004	7.7E-003 C	
ACETONE	67641	1.00E-001 I				y	6.1E+002 N	3.7E+002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N	1.2E-001	2.5E+000 N	
ACETONITRILE	75058			1.7E-002 I		y	1.2E+002 N	6.2E+001 N						
ACETOPHENONE	98862	1.00E-001 I		5.70E-006 W		y	4.2E-002 N	2.1E-002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N	2.9E-002	5.8E-001 N	
ACROLEIN	107028	2.00E-002 H		5.70E-006 I		y	4.2E-002 N	1.5E-002 C	2.7E+001 N	4.1E+004 N	1.6E+003 N	1.1E-005	2.2E-004 N	
ACRYLAMIDE	79061	2.00E-004 I	4.50E+000 I		4.50E+000 I	y	8.4E-001 C	1.4E-002 C	7.0E-004 C	1.3E+000 C	1.4E-001 C	1.0E-005	2.0E-004 N	
ACRYLONITRILE	107131	1.00E-003 H	5.40E-001 I	5.70E-004 I	2.40E-001 I	y	3.7E-002 C	2.6E-002 C	5.8E-003 C	1.1E+001 C	1.2E+000 C	3.7E-006	7.4E-005 C	
ALACHLOR	15972608	1.00E-002 I	8.00E-002 H				5.5E+003 N	5.5E+002 N	2.0E+002 N	3.1E+005 N	1.2E+004 N			
ALAR	1596845	1.50E-001 I					3.7E+001 N	3.7E+000 N	1.4E+000 N	2.0E+003 N	7.8E+001 N	1.0E-002	2.1E-001 N	
ALDICARB	116063	1.00E-003 I					3.7E+001 N	3.7E+000 N	1.4E+000 N	2.0E+003 N	7.8E+001 N			
ALDICARB SULFONE	1646884	1.00E-003 I					3.9E-003 C	3.7E-004 C	1.9E-004 C	2.0E+006 N	7.8E+004 N	3.8E-004	7.7E-003 C	
ALDRIN	309002	3.00E-005 I	1.70E+001 I		1.70E+001 I		7.3E-001 N	7.3E-002 N	8.1E-002 N	1.2E+002 N	4.7E+000 N			
ALUMINUM	7429905	1.00E+000 E		1.00E-003 E			2.1E+002 N	1.0E+002 N	1.2E+001 C	3.4E-001 C	3.8E-002 C			
AMINODINITROTOLUENES		6.00E-005 E					3.7E+004 N	3.7E+000 N	1.4E+003 N	2.0E+006 N	7.8E+004 N			
4-AMINOPYRIDINE	504245	2.00E-005 H					2.2E+000 N	2.2E-001 N	8.1E-002 N	1.2E+002 N	4.7E+000 N			
AMMONIA	7664417			2.86E-002 I		y	2.1E+002 N	1.0E+002 N	1.2E+001 C	3.4E-001 C	3.8E-002 C			
ANILINE	62533	7.00E-003 E	5.70E-003 I	2.90E-004 I			1.2E+001 C	1.1E+000 N	5.5E-001 C	1.0E+003 C	1.1E+002 C I	6.8E-003	1.4E-001 C	
ANTIMONY	7440360	4.00E-004 I					1.5E+001 N	1.5E+000 N	5.4E-001 N	8.2E+002 N	3.1E+001 N	6.6E-001	1.3E+001 N	
ANTIMONY PENTOXIDE	1314609	5.00E-004 H					1.8E+001 N	1.8E+000 N	6.8E-001 N	1.0E+003 N	3.9E+001 N			
ANTIMONY TETROXIDE	1332816	4.00E-004 H					1.5E+001 N	1.5E+000 N	5.4E-001 N	8.2E+002 N	3.1E+001 N			
ANTIMONY TRIOXIDE	1309644	4.00E-004 H		5.70E-005 I			1.5E+001 N	2.1E-001 N	5.4E-001 N	8.2E+002 N	3.1E+001 N			
ARSENIC	7440382	3.00E-004 I	1.50E+000 I		1.51E+001 I		4.5E-002 C	4.1E-004 C	2.1E-003 C	3.8E+000 C	4.3E-001 C	1.3E-003	2.6E-002 C	
ARSINE	7784421			1.40E-005 I		y	1.0E-001 N	5.1E-002 N	1.2E+001 N	3.1E+001 N				
ASSURE	76578148	9.00E-003 I					3.3E+002 N	3.3E+001 N	1.2E+001 N	1.8E+004 N	7.0E+002 N			
ATRAZINE	1912249	3.50E-002 I	2.20E-001 H				3.0E-001 C	2.8E-002 C	1.4E-002 C	2.6E+001 C	2.9E+000 C	4.4E-004	8.8E-003 C	
AZOBENZENE	103333			1.10E-001 I			6.1E-001 C	5.7E-002 C	2.9E-002 C	5.2E+001 C	5.8E+000 C	1.8E-003	3.5E-002 C	
BARIUM	7440393	7.00E-002 I			1.10E-001 I		2.6E+003 N	5.1E-001 N	9.5E+001 N	1.4E+005 N	5.5E+003 N	1.1E+002	2.1E+003 N	
BAYGON	114261	4.00E-003 I					1.5E+002 N	1.5E+001 N	5.4E+000 N	8.2E+003 N	3.1E+002 N			
BAYTHROID	68359375	2.50E-002 I					9.1E+002 N	9.1E+001 N	3.4E+001 N	5.1E+004 N	2.0E+003 N			
BENTAZON	25057890	3.00E-002 I					1.1E+001 N	1.1E+002 N	4.1E+001 N	6.1E+004 N	2.3E+003 N			
BENZALDEHYDE	100527	1.00E-001 I					3.7E+003 N	3.7E+002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N			
BENZENE	71432	3.00E-003 E	5.5E-002 I	1.70E-003 E	2.90E-002 I	y	6.1E-002 N	3.2E-001 C	2.2E-001 C	5.7E-002 C	1.0E+002 C	1.2E+001 C	9.0E-005	1.8E-003 C
BENZENETHIOL	108985	1.00E-005 H					2.9E-004 C	1.5E+005 N	1.5E+004 N	5.4E+003 N	8.2E+006 N	3.1E+005 N		
BENZIDINE	92875	3.00E-003 I	2.30E+002 I		2.30E+002 I		1.5E+005 N	1.5E+004 N	5.4E+003 N	8.2E+006 N	3.1E+005 N			
BENZOIC ACID	65850	4.00E+000 I					1.1E+004 N	1.1E+003 N	4.1E+002 C	6.1E+005 N	2.3E+004 N	4.4E+000	8.8E+001 N	
BENZYL ALCOHOL	100516	3.00E-001 H					6.2E-002 C	3.7E-002 N	1.4E-002 N	2.0E+001 N	7.8E-001 N			
BENZYL CHLORIDE	100447		0.17 I				7.3E-001 N	7.5E-004 C	2.7E+000 N	4.1E+001 C	4.1E+002 C	1.9E-005	3.7E-004 C	
BERYLLIUM	7440417	2.00E-003 I		5.7E-006 I	8.40E+000 I	y	3.0E+002 N	3.7E-002 C	1.9E-002 C	3.4E+001 C	3.8E+000 C	1.9E-005	3.7E-004 C	
BIPHENYL	92524	5.00E-002 I					1.8E+002 N	1.8E+001 N	7.5E-004 C	2.7E+000 N	4.1E+003 N	1.6E+002 N	5.8E+001	1.2E+003 N
BIS(2-CHLOROETHYL)ETHER	111444		1.10E+000 I		1.10E+000 I	y	3.0E+002 N	1.8E+002 N	6.8E+001 N	1.0E+005 N	3.9E+003 N	4.8E+000	9.6E+001 N	
BIS(2-CHLOROISOPROPYL)ETHER	108601	4.00E-002 I	7.00E-002 H		3.50E-002 H	y	2.8E-001 C	1.8E-001 C	2.9E-003 C	5.2E+000 C	5.8E-001 C	2.2E-006	4.4E-005 C	
BIS(CHLOROMETHYL)ETHER	542881		2.20E+002 I		2.20E+002 I	y	4.8E-005 C	2.8E-005 C	4.5E-002 C	8.2E+001 C	9.1E+000 C	8.4E-005	1.7E-003 C	
BIS(2-ETHYLHEXYL)PHthalate	117817	2.00E-002 I	1.40E-002 I		1.40E-002 E		4.8E+000 C	4.5E-001 C	2.3E-001 C	4.1E+002 C	4.6E+001 C	9.7E-009	1.9E-007 C	
BORON	7440428	9.00E-002 I			5.70E-003 H		3.3E+003 N	2.1E+001 N	1.2E+002 N	1.8E+005 N	7.0E+003 N	1.4E+002	2.9E+003 C	

Sources: I = IRIS; H = HEAST; A = HEAST Alternate; W = Withdrawn from IRIS or HEAST E = EPA-NCEA provisional value; O = other												Basis: C = Carcinogenic effects; N = Noncarcinogenic effects; I = RBC at HI of 0.1 < RBC < 1		Region III SSLs									
Chemical	CAS	RfDo		CSFo		RfDi		CSFi		VOC		Risk-based concentrations											
		mg/kg/d	1/mg/kg/d	mg/kg/d	1/mg/kg/d	mg/kg/d	1/mg/kg/d	ug/l	Tap water	Ambient air ug/m3	Fish mg/kg	Soil Industrial mg/kg	Residential mg/kg	DAF 1 mg/kg	DAF 20 mg/kg								
BROMODICHLOROMETHANE	75274	2.00E-002	I	6.20E-002	I			y	1.7E-001	C	1.0E-001	C	5.1E-002	C	9.2E+001	C	1.0E+001	C	5.4E-005	1.1E-003	C		
BROMOETHENE	593602					8.6E-004	I	1.10E-001	H	y	1.1E-001	C	5.7E-002	C	4.0E-001	C	7.2E+002	C	8.1E+001	C	5.4E-005	1.1E-003	C
BROMOFORM	75252	2.00E-002	I	7.90E-003	I			3.90E-003	I		8.5E+000	C	1.6E+000	C	1.9E+000	N	2.9E+003	N	1.1E+002	N	3.3E-003	6.7E-002	C
BROMOMETHANE	74839	1.40E-003	I			1.40E-003	I			y	8.5E+000	N	5.1E+000	N	1.9E+000	N	1.0E+004	N	3.9E+002	N	2.1E-003	4.1E-002	N
BROMOPHOS	2104963	5.00E-003	H								1.8E+002	N	1.8E+001	N	6.8E+000	N					3.9E-006	7.8E-005	C
1,3-BUTADIENE	106990								1.80E+000	H	y	7.0E-003	C	3.5E-003	C								
1-BUTANOL	71363	1.00E-001	I								3.7E+003	N	3.7E+002	N	1.4E+002	N	2.0E+005	N	7.8E+003	N	7.8E-001	1.6E+001	N
BUTYLBENZYLPHthalATE	85687	2.00E-001	I								7.3E+003	N	7.3E+002	N	2.7E+002	N	4.1E+005	N	1.6E+004	N	8.4E+002	1.7E+004	N
BUTYLATE	2008415	5.00E-002	I								1.8E+003	N	1.8E+002	N	6.8E+001	N	1.0E+005	N	3.9E+003	N			
N-BUTYLBENZENE	104518	4.00E-002	E							y	2.4E+002	N	1.5E+002	N	5.4E+001	N	8.2E+004	N	3.1E+003	N			
SEC-BUTYLBENZENE	135988	4.00E-002	E							y	2.4E+002	N	1.5E+002	N	5.4E+001	N	8.2E+004	N	3.1E+003	N			
TERT-BUTYLBENZENE	98066	4.00E-002	E							y	2.4E+002	N	1.5E+002	N	5.4E+001	N	8.2E+004	N	3.1E+003	N			
CADMIUM-WATER	7440439	5.00E-004	I			5.7E-005	E	6.30E+000	I		1.8E+001	N	9.9E-004	C	6.8E-001	N	1.0E+003	N	3.9E+001	N	1.4E+000	2.7E+001	N
CADMIUM-FOOD	7440439	1.00E-003	I			5.7E-005	E	6.30E+000	I		3.7E+001	N	9.9E-004	C	1.4E+000	N	2.0E+003	N	7.8E+001	N	2.7E+000	5.5E+001	N
CAPROLACTAM	105602	5.00E-001	I								1.8E+004	N	1.8E+003	N	6.8E+002	N	1.0E+006	N	3.9E+004	N			
CARBARYL	63252	1.00E-001	I								3.7E+003	N	3.7E+002	N	1.4E+002	N	2.0E+005	N	7.8E+003	N	1.5E+000	3.0E+001	N
CARBON DISULFIDE	75150	1.00E-001	I			2.00E-001	I			y	1.0E+003	N	7.3E+002	N	1.4E+002	N	2.0E+005	N	7.8E+003	N	9.5E-001	1.9E+001	N
CARBON TETRACHLORIDE	56235	7.00E-004	I	1.30E-001	I	5.71E-004	E	5.30E-002	I	y	1.6E-001	C	1.2E-001	C	2.4E-002	C	4.4E+001	C	4.9E+000	C	1.1E-004	2.1E-003	C
CARBOSULFAN	55285148	1.00E-002	I								3.7E+002	N	3.7E+001	N	1.4E+002	N	2.0E+005	N	7.8E+003	N			
CHLORAL HYDRATE	302170	1.00E-001	I								3.7E+003	N	3.7E+002	N	1.4E+002	N	2.0E+005	N	7.8E+003	N			
CHLORANIL	118752			4.00E-001	H						1.7E-001	C	1.6E-002	C	7.9E-003	C	1.4E+001	C	1.6E+000	C			
CHLORDANE	57749	5.00E-004	I	3.5E-001	I	2.00E-004	I	3.5E-001	I		1.9E-001	C	1.8E-002	C	9.0E-003	C	1.6E+001	C	1.8E+000	C	4.6E-002	9.2E-001	C
CHLORINE	7782505	1.00E-001	I			5.7E-005	E			y	4.2E+001	N	2.1E+001	N	1.4E+002	N	2.0E+005	N	7.8E+003	N			
**CHLORINE DIOXIDE	1040944	3.00E-002	I			5.70E-005	I			y	4.2E+001	N	2.1E+001	N	4.1E+001	N	6.1E+004	N	2.3E+003	N			
CHLOROACETIC ACID	79118	2.00E-003	H								7.3E+001	N	7.3E+000	N	2.7E+000	N	4.1E+003	N	1.6E+002	N			
4-CHLOROANILINE	106478	4.00E-003	I								1.5E+002	N	1.5E+001	N	5.4E+000	N	8.2E+003	N	3.1E+002	N	4.8E-002	9.7E-001	N
CHLOROBENZENE	108907	2.00E-002	I			1.7E-002	E			y	1.1E+002	N	6.2E+001	N	2.7E+001	N	4.1E+004	N	1.6E+003	N	4.0E-002	8.0E-001	N
CHLOROBENZILATE	510156	2.00E-002	I	2.70E-001	H			2.70E-001	H		2.5E-001	C	2.3E-002	C	1.2E-002	C	2.1E+001	C	2.4E+000	C	1.3E-003	2.7E-002	C
P-CHLOROBENZOIC ACID	74113	2.00E-001	H								7.3E+003	N	7.3E+002	N	2.7E+002	N	4.1E+005	N	1.6E+004	N			
2-CHLORO-1,3-BUTADIENE	126998	2.00E-002	A			2.00E-003	H			y	1.4E+001	N	7.3E+000	N	2.7E+001	N	4.1E+004	N	1.6E+003	N	6.0E-003	1.2E-001	N
1-CHLOROBUTANE	109693	4.00E-001	H								2.4E+003	N	1.5E+003	N	5.4E+002	N	8.2E+005	N	3.1E+004	N	1.0E+000	2.0E+001	N
1-CHLORO-1,1-DIFLUOROETHANE	75683					1.40E+001	I			y	1.0E+005	N	5.1E+004	N							7.0E+001	1.4E+003	N
CHLORODIFLUOROMETHANE	75456					1.40E+001	I			y	1.0E+005	N	5.1E+004	N							7.0E+001	1.4E+003	N
CHLOROETHANE	75003	4.00E-001	E	2.90E-003	E	2.90E+000	I			y	3.6E+000	C	2.2E+000	C	1.1E+000	C	2.0E+003	C	2.2E+002	C	9.6E-004	1.9E-002	C
CHLOROFORM	67663	1.00E-002	I	6.10E-003	I	8.6E-005	E	8.10E-002	I	y	1.5E-001	C	7.7E-002	C	5.2E-001	C	9.4E+002	C	1.0E+002	C	4.5E-005	8.9E-004	C
CHLORMETHANE	74873			1.30E-002	H	8.6E-002	E	3.5E-003	E	y	2.1E+000	C	1.8E+000	C	2.4E-001	C	4.4E+002	C	4.9E+001	C	5.2E-004	1.0E-002	C
4-CHLORO-2-METHYLANILINE	95692			5.80E-001	H						1.2E-001	C	1.1E-002	C	5.4E-003	C	9.9E+000	C	1.1E+000	C			
BETA-CHLORONAPHTHALENE	91587	8.00E-002	I								4.9E+002	N	2.9E+002	N	1.1E+002	N	1.6E+005	N	6.3E+003	N	1.6E+000	3.2E+001	N
O-CHLORONITROBENZENE	88733			2.50E-002	H						4.2E-001	C	2.5E-001	C	1.3E-001	C	2.3E+002	C	2.6E+001	C			
P-CHLORONITROBENZENE	100005					1.80E-002	H				5.9E-001	C	3.5E-001	C	1.8E-001	C	3.2E+002	C	3.5E+001	C			
2-CHLOROPHENOL	95578	5.00E-003	I								3.0E+001	N	1.8E+001	N	6.8E+000	N	1.0E+004	N	3.9E+002	N			
2-CHLOROPROPANE	75296					2.90E-002	H				2.1E+002	N	1.1E+002	N							6.6E-002	1.3E+000	N
O-CHLORTOLUENE	95498	2.00E-002	I								1.2E+002	N	7.3E+001	N	2.7E+001	N	4.1E+004	N	1.6E+003	N	6.5E-002	1.3E+000	N
CHLORPYRIFOS	2921882	3.00E-003	I								1.1E+002	N	1.1E+001	N	4.1E+000	N	6.1E+003	N	2.3E+002	N	3.2E+000	6.3E+001	N
CHLORPYRIFOS-METHYL	5598130	1.00E-002	H								3.7E+002	N	3.7E+001	N	1.4E+001	N	2.0E+004	N	7.8E+002	N			

Sources: I = IRIS H = HEAST A = HEAST Alternate W = Withdrawn from IRIS or HEAST  
 E = EPA-NCEA provisional value O = other

Chemical	CAS	RfDo	CSFo	RfDi	CSFi	VOC	Tap water ug/l	Ambient air ug/m <sup>3</sup>	Risk-based concentrations			Region III SSS	
		mg/kg/d	1/mg/kg/d	mg/kg/d	1/mg/kg/d				Fish mg/kg	Soil Industrial mg/kg	Residential mg/kg	DAF 1 mg/kg	DAF 20 mg/kg
CHROMIUM III	16065831	1.50E+000 I					5.5E+004 N	5.5E+003 N	2.0E+003 N	3.1E+006 N	1.2E+005 N	9.9E+007	2.0E+009 N
CHROMIUM VI	18540299	3.00E-003 I		3.00E-005 I	4.10E+001 H		1.1E+002 N	1.5E-004 C	4.1E+000 N	6.1E+003 N	2.3E+002 N	2.1E+000	4.2E+001 N
**COBALT	7440484	2.00E-002 E		5.00E-006 E			7.3E+002 N	1.8E-002 N	2.7E+001 N	4.1E+004 N	1.6E+003 N		
COKE OVEN EMISSIONS (COAL TAR)	8007452				2.2 I		2.8E-003 C						
COPPER	7440508	4.00E-002 H					1.5E+003 N	1.5E+002 N	5.4E+001 N	8.2E+004 N	3.1E+003 N	5.3E+002	1.1E+004 N
CROTONALDEHYDE	123739		1.90E+000 H				5.6E-003 C	3.3E-003 C	1.7E-003 C	3.0E+000 C	3.4E-001 C	1.5E-005	3.1E-004 C
CUMENE	98828	1.00E-001 I		1.10E-001 I			6.6E-002 N	4.0E+002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N	3.2E+000	6.4E+001 N
CYANIDE (FREE)	57125	2.00E-002 I					7.3E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N		
CALCIUM CYANIDE	592018	4E-002 I					1.5E+003 N	1.5E+002 N	5.4E+001 N	8.2E+004 N	3.1E+003 N		
COPPER CYANIDE	544923	5.00E-003 I					1.8E+002 N	1.8E+001 N	6.8E+000 N	1.0E+004 N	3.9E+002 N		
CYANAZINE	21725462	2.00E-003 H	8.40E-001 H				8.0E-002 C	7.5E-003 C	3.8E-003 C	6.8E+000 C	7.6E-001 C	2.6E-005	5.3E-004 C
CYANOGEN	460195	4.00E-002 I					2.4E+002 N	1.5E+002 N	5.4E+001 N	8.2E+004 N	3.1E+003 N		
CYANOGEN BROMIDE	506683	9.00E-002 I					3.3E+003 N	3.3E+002 N	1.2E+002 N	1.8E+005 N	7.0E+003 N		
CYANOGEN CHLORIDE	506774	5.00E-002 I					1.8E+003 N	1.8E+002 N	6.8E+001 N	1.0E+005 N	3.9E+003 N		
HYDROGEN CYANIDE	74908	2.00E-002 I		8.60E-004 I			6.2E+000 N	3.1E+000 N	2.7E+001 N	4.1E+004 N	1.6E+003 N	1.1E-001	2.2E+000 N
POTASSIUM CYANIDE	151508	5.00E-002 I					1.8E+003 N	1.8E+002 N	6.8E+001 N	1.0E+005 N	3.9E+003 N		
POTASSIUM SILVER CYANIDE	506616	2.00E-001 I					1.8E+003 N	1.8E+002 N	6.8E+001 N	1.0E+005 N	3.9E+003 N		
SILVER CYANIDE	506649	1.00E-001 I					7.3E+003 N	7.3E+002 N	2.7E+002 N	4.1E+005 N	1.6E+004 N		
SODIUM CYANIDE	143339	4.00E-002 I					3.7E+003 N	3.7E+002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N	3.1E+001	6.2E+002 N
THIOCYANATE		5.00E-002 E					1.5E+003 N	1.5E+002 N	5.4E+001 N	8.2E+004 N	3.1E+003 N		
ZINC CYANIDE	557211	5.00E-002 I					1.8E+003 N	1.8E+002 N	6.8E+001 N	1.0E+005 N	3.9E+003 N		
CYCLOHEXANONE	108941	5.00E+000 I					1.8E+003 N	1.8E+002 N	6.8E+001 N	1.0E+005 N	3.9E+003 N	1.1E+002	2.3E+003 N
CYHALOTHIN/KARATE	68085858	5.00E-003 I					1.8E+005 N	1.8E+004 N	6.8E+003 N	1.0E+007 N	3.9E+005 N	6.1E+001	1.2E+003 N
CYPERMETHRIN	52315078	1.00E-002 I					1.8E+002 N	1.8E+001 N	6.8E+000 N	1.0E+004 N	3.9E+002 N		
DACTHAL	1861321	1.00E-002 I					3.7E+002 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N		
DALAPON	75990	3.00E-002 I					3.7E+002 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N		
DDD	72548		2.40E-001 I				1.1E+003 N	1.1E+002 N	4.1E+001 N	6.1E+004 N	2.3E+003 N	3.5E-001	7.1E+000 N
DDOE	72559		3.40E-001 I				2.8E-001 C	2.6E-002 C	1.3E-002 C	2.4E+001 C	2.7E+000 C	5.6E-001	1.1E+001 C
DDT	50293	5.00E-004 I	3.40E-001 I		3.40E-001 I		2.0E-001 C	1.8E-002 C	9.3E-003 C	1.7E+001 C	1.9E+000 C	1.8E+000	3.5E+001 C
DAZINON	333415	9.00E-004 H					2.0E-001 C	1.8E-002 C	9.3E-003 C	1.7E+001 C	1.9E+000 C	5.8E-002	1.2E+000 C
DIBENZOFURAN	132649	4.00E-003 E					3.3E+001 N	3.3E+000 N	1.2E+000 N	1.8E+003 N	7.0E+001 N	2.1E-002	4.3E-001 N
1,4-DIBROMOBENZENE	106376	1.00E-002 I					2.4E+001 N	1.5E+001 N	5.4E+000 N	8.2E+003 N	3.1E+002 N	3.8E-001	7.7E+000 N
DIBROMOCHLOROMETHANE	124481	2.00E-002 I	8.40E-002 I				3.7E+002 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N		
1,2-DIBROMO-3-CHLOROPROPANE	96128		1.40E+000 H	5.70E-005 I	2.40E-003 H	y	1.3E-001 C	7.5E-002 C	3.8E-002 C	6.8E+001 C	7.6E+000 C	4.1E-005	8.3E-004 C
1,2-DIBROMOETHANE	106934		8.50E+001 I	5.70E-005 H	7.60E-001 I	y	4.7E-002 C	2.1E-001 N	2.3E-003 C	4.1E+000 C	4.6E-001 C	4.4E-005	8.7E-004 C
DI BUTYL PHTHALATE	84742	1.00E-001 I					7.5E-004 C	8.2E-003 C	3.7E-005 C	6.7E-002 C	7.5E-003 C	4.3E-007	8.5E-006 C
DICAMBA	1918009	3.00E-002 I					3.7E+003 N	3.7E+002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N	2.5E+002	5.0E+003 N
1,2-DICHLOROBENZENE	95501	9.00E-002 I					1.1E+003 N	1.1E+002 N	4.1E+001 N	6.1E+004 N	2.3E+003 N	2.2E-001	4.5E+000 N
**1,3-DICHLOROBENZENE	541731	3.00E-002 E					5.5E+002 N	3.3E+002 N	1.2E+002 N	1.8E+005 N	7.0E+003 N	4.6E-001	9.3E+000 N
1,4-DICHLOROBENZENE	106467	3.00E-002 E	2.40E-002 H	2.29E-001 I	2.2E-002 E	y	1.8E+002 N	1.1E+002 N	4.1E+001 N	6.1E+004 N	2.3E+003 N	1.5E-001	2.9E+000 N
3,3'-DICHLOROBENZIDINE	91941		4.50E-001 I				4.7E-001 C	2.8E-001 C	1.3E-001 C	2.4E+002 C	2.7E+001 C	3.6E-004	7.1E-003 C
1,4-DICHLORO-2-BUTENE	764410						1.5E-001 C	1.4E-002 C	7.0E-003 C	1.3E+001 C	1.4E+000 C	2.5E-004	4.9E-003 C
DICHLORODIFLUOROMETHANE	75718	2.00E-001 I		5.00E-002 A		y	1.3E-003 C	6.7E-004 C				4.0E-007	8.0E-006 C
1,1-DICHLOROETHANE	75343	1.00E-001 H		1.40E-001 A		y	3.5E+002 N	1.8E+002 N	2.7E+002 N	4.1E+005 N	1.8E+004 N	5.5E-001	1.1E+001 N
1,2-DICHLOROETHANE	107062	3.00E-002 E	9.10E-002 I	1.40E-003 E	9.10E-002 I	y	8.0E+002 N	5.1E+002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N	2.3E-001	4.5E+000 N
							1.2E-001 C	6.9E-002 C	3.5E-002 C	6.3E+001 C	7.0E+000 C	5.2E-005	1.0E-003 C

Sources: I = IRIS H = HEAST A = HEAST Alternate W = Withdrawn from IRIS or HEAST E = EPA-NCEA provisional value O = other										Basis: C = Carcinogenic effects N = Noncarcinogenic effects I = RBC at HI of 0.1 < RBC-C Risk-based concentrations					Region III SSLs Soil, for groundwater migration	
Chemical	CAS	RfDo	CSFo	RfDi	CSFi	VOC	Tap water ug/l	Ambient air ug/m3	Fish mg/kg	Soil Industrial mg/kg	Residential mg/kg	DAF 1 mg/kg	DAF 20 mg/kg			
		mg/kg/d	1/mg/kg/d	mg/kg/d	1/mg/kg/d											
1,1-DICHLOROETHENE	75354	9.00E-003 I	6.00E-001 I		1.75E-001 I	y	4.4E-002 C	3.6E-002 C	5.3E-003 C	9.5E+000 C	1.1E+000 C	1.8E-005	3.6E-004 C			
CIS-1,2-DICHLOROETHENE	156592	1.00E-002 H				y	6.1E+001 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N	1.7E-002	3.5E-001 N			
TRANS-1,2-DICHLOROETHENE	156805	2.00E-002 I				y	1.2E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N	4.1E-002	8.2E-001 N			
TOTAL 1,2-DICHLOROETHENE	540590	9.00E-003 H				y	5.5E+001 N	3.3E+001 N	1.2E+001 N	1.8E+004 N	7.0E+002 N	1.9E-002	3.7E-001 N			
2,4-DICLOROPHENOL	120832	3.00E-003 I					1.1E+002 N	1.1E+001 N	4.1E+000 N	6.1E+003 N	2.3E+002 N	6.0E-002	1.2E+000 N			
2,4-D	94757	1.00E-002 I					3.7E+002 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N	4.5E-001	9.0E+000 N			
4-(2,4-DICHLOROPHOXY)BUTYRIC ACID	94826	8E-003 I					2.9E+002 N	2.9E+001 N	1.1E+001 N	1.6E+004 N	6.3E+002 N					
1,2-DICHLOROPROPANE	78875		6.80E-002 H	1.14E-003 I		y	1.6E-001 C	9.2E-002 C	4.6E-002 C	8.4E-001 C	9.4E+000 C	1.0E-004	2.1E-003 C			
2,3-DICHLOROPROPANOL	616239	3.00E-003 I					1.1E+002 N	1.1E+001 N	4.1E+000 N	6.1E+003 N	2.3E+002 N					
1,3-DICHLOROPROPENE	542756	3.00E-002 I	1.00E-001 I	5.71E-003 I	1.00E-002 I	y	4.4E-001 C	6.3E-001 C	3.2E-002 C	5.7E+001 C	6.4E+000 C	1.6E-004	3.1E-003 C			
DICHLORVOS	62737	5E-004 I	0.29 I	1.43E-004 I			2.3E-001 C	2.2E-002 C	1.1E-002 C	2.0E+001 C	2.2E+000 C	5.5E-005	1.1E-003 C			
DICOFOL	115322		4.4E-001 W				1.5E-001 C	1.4E-002 C	7.2E-003 C	1.3E+001 C	1.5E+000 C	9.3E-004	1.9E-002 C			
DICYCLOPENTADIENE	77736	3E-002 H		6.00E-005 A		y	4.4E-001 N	2.2E-001 N	4.1E+001 N	6.1E+004 N	2.3E+003 N					
DIELDRIN	60571	5.00E-005 I	1.60E+001 I		1.60E+001 I		4.2E-003 C	3.9E-004 C	2.0E-004 C	3.6E-001 C	4.0E-002 C	1.1E-004	2.2E-003 C			
DIESEL EMISSIONS				1.40E-003 I			5.1E+000 N									
DIETHYLPHthalate	84662	8.00E-001 I					2.9E+004 N	2.9E+003 N	1.1E+003 N	1.6E+006 N	6.3E+004 N	2.3E+001	4.5E+002 N			
DIETHYLENE GLYCOL, MONOBUTYL ETHER	112345			5.70E-003 H			2.1E+001 N									
DIETHYLENE GLYCOL, MONOETHYL ETHER	111900	2.00E+000 H					7.3E+004 N	7.3E+003 N	2.7E+003 N	4.1E+006 N	1.6E+005 N					
DI(2-ETHYLHEXYL)ADIPATE	103231	6.00E-001 I	1.20E-003 I				5.6E+001 C	5.2E+000 C	2.6E+000 C	4.8E+003 C	5.3E+002 C					
DIETHYLSTILBESTROL	56531		4.70E+003 H				1.4E-005 C	1.3E-006 C	6.7E-007 C	1.2E-003 C	1.4E-004 C					
DIFENOZOQUAT (AVENGE)	43222486	8.00E-002 I					2.9E+002 N	1.1E+002 N	1.6E+005 N	6.3E+003 N						
1,1-DIFLUOROETHANE	75376		1.10E+001 I			y	8.0E+004 N	4.0E+004 N								
DIISOPROPYL Methylphosphonate (DIMP)	1445758	8.00E-002 I					2.9E+003 N	2.9E+002 N	1.1E+002 N	1.6E+005 N	6.3E+003 N					
3,3'-DIMETHOXYBENZIDINE	119904		1.40E-002 H		5.70E-006 W	y	4.8E+000 C	4.5E-001 C	2.3E-001 C	4.1E+002 C	4.6E+001 C		8.5E-006	1.7E-004 N		
DIMETHYLAMINE	124403						4.2E-002 N	2.1E-002 N								
2,4-DIMETHYLANILINE HYDROCHLORIDE	21436964		5.80E-001 H				1.2E-001 C	1.1E-002 C	5.4E-003 C	9.9E+000 C	1.1E+000 C					
2,4-DIMETHYLANILINE	95681		7.50E-001 H				8.9E-002 C	8.3E-003 C	4.2E-003 C	7.6E+000 C	8.5E-001 C					
N,N-DIMETHYLANILINE	121697	2.00E-003 I					7.3E+001 N	7.3E+000 N	2.7E+000 N	4.1E+003 N	1.6E+002 N					
3,3'-DIMETHYLBENZIDINE	119937		9.20E+000 H				7.3E-003 C	6.8E-004 C	3.4E-004 C	6.2E-001 C	6.9E-002 C					
1,1-DIMETHYLHYDRAZINE	57147		2.60E+000 W		3.50E+000 W		2.6E-002 C	1.8E-003 C	1.2E-003 C	2.2E+000 C	2.5E-001 C					
1,2-DIMETHYLHYDRAZINE	540738		3.70E+001 W		3.70E+001 W		1.8E-003 C	1.7E-004 C	8.5E-005 C	1.5E-001 C	1.7E-002 C					
2,4-DIMETHYLPHENOL	105679	2.00E-002 I					7.3E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N		3.4E-001	6.7E+000 N		
2,6-DIMETHYLPHENOL	576261	6.00E-004 I					2.2E+001 N	2.2E+000 N	8.1E-001 N	1.2E+003 N	4.7E+001 N					
3,4-DIMETHYLPHENOL	95658	1.00E-003 I					3.7E+001 N	3.7E+000 N	1.4E+000 N	2.0E+003 N	7.8E+001 N					
DIMETHYLPHthalate	131113	1.00E+001 W					3.7E+005 N	3.7E+004 N	1.4E+004 N	2.0E+007 N	7.8E+005 N					
1,2-DINITROBENZENE	528290	4.00E-004 H					1.5E+001 N	1.5E+000 N	5.4E-001 N	8.2E+002 N	3.1E+001 N					
1,3-DINITROBENZENE	99650	1.00E-004 I					3.7E+000 N	3.7E-001 N	1.4E-001 N	2.0E+002 N	7.8E+000 N		1.8E-003	3.7E-002 N		
1,4-DINITROBENZENE	100254	4.00E-004 H					1.5E+001 N	1.5E+000 N	5.4E-001 N	8.2E+002 N	3.1E+001 N					
4,6-DINITRO-O-CYCLOHEXYL PHENOL	131895	2.00E-003 I					7.3E+001 N	7.3E+000 N	2.7E+000 N	4.1E+003 N	1.6E+002 N					
4,6-DINITRO-2-METHYLPHENOL	534521	1.00E-004 E					3.7E+000 N	3.7E-001 N	1.4E-001 N	2.0E+002 N	7.8E+000 N					
2,4-DINITROPHENOL	51285	2.00E-003 I			6.80E-001 I		7.3E+001 N	7.3E+000 N	2.7E+000 N	4.1E+003 N	1.6E+002 N					
DINITROTOLUENE MIX							9.8E-002 C	9.2E-003 C	4.6E-003 C	8.4E+000 C	9.4E-001 C					
2,4-DINITROTOLUENE	121142	2.00E-003 I					7.3E+001 N	7.3E+000 N	2.7E+000 N	4.1E+003 N	1.6E+002 N	2.9E-002	5.7E-001 N			
2,6-DINITROTOLUENE	606202	1.00E-003 H					3.7E+001 N	3.7E+000 N	1.4E+000 N	2.0E+003 N	7.8E+001 N	1.2E-002	2.5E-001 N			
DINOSEB	88857	1.00E-003 I					3.7E+001 N	3.7E+000 N	1.4E+000 N	2.0E+003 N	7.8E+001 N	8.7E-003	1.7E-001 N			

Sources I = IRIS H = HEAST A = HEAST Alternate W = Withdrawn from IRIS or HEAST  
 E = EPA-NCEA provisional value O = other

Chemical	CAS	RfDo mg/kg/d	CSFo 1/mg/kg/d	RfDi mg/kg/d	CSFi 1/mg/kg/d	VOC	Risk-based concentrations					Region III SSLs	
							Tap water ug/l	Ambient air ug/m3	Fish mg/kg	Soil Industrial mg/kg	Residential mg/kg	Soil, for groundwater migration DAF 1 mg/kg	DAF 20 mg/kg
DIOCYLPHthalATE	117840	2.00E-002 H					7.3E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N	1.2E+005	2.4E+006 N
1,4-DIOXANE	123911		1.10E-002 I				6.1E+000 C	5.7E-001 C	2.9E-001 C	5.2E+002 C	5.8E+001 C	1.3E-003	2.6E-002 C
DIPHENYLAMINE	122394	2.50E-002 I					9.1E+002 N	9.1E+001 N	3.4E+001 N	5.1E+004 N	2.0E+003 N	1.3E+000	2.5E+001 N
1,2-DIPHENYLHYDRAZINE	122667			8.00E-001 I			8.4E-002 C	7.8E-003 C	3.9E-003 C	7.2E+000 C	8.0E-001 C	1.3E-004	2.5E-003 C
DIQUAT	85007	2.20E-003 I					8.0E+001 N	8.0E+000 N	3.0E+000 N	4.5E+003 N	1.7E+002 N	1.7E-002	3.3E-001 N
DISULFOTON	298044	4.00E-005 I					1.5E+000 N	1.5E-001 N	5.4E-002 N	8.2E+001 N	3.1E+000 N	3.2E-003	6.4E-002 N
1,4-DITHIANE	505293	1.00E-002 I					3.7E+002 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N		
DIURON	330541	2.00E-003 I					7.3E+001 N	7.3E+000 N	2.7E+000 N	4.1E+003 N	1.6E+002 N	5.8E-002	1.2E+000 N
ENDOSULFAN	115297	6.00E-003 I					2.2E+002 N	2.2E+001 N	8.1E+000 N	1.2E+004 N	4.7E+002 N	9.8E-001	2.0E+001 N
ENDRIN	72208	3.00E-004 I					1.1E+001 N	1.1E+000 N	4.1E-001 N	6.1E+002 N	2.3E+001 N	2.7E-001	5.4E+000 N
EPICHLOROHYDRIN	106898	2.00E-003 H	9.90E-003 I	2.86E-004 I	4.20E-003 I	y	2.0E+000 N	1.0E+000 N	3.2E-001 C	5.8E+002 C	6.5E+001 C	4.2E-004	8.4E-003 N
ETHION	563122	5.00E-004 I					1.8E+001 N	1.8E+000 N	6.8E-001 N	1.0E+003 N	3.9E+001 N	3.2E-001	6.4E+000 N
2-ETHOXYETHANOL	110805	4.00E-001 H		5.70E-002 I			1.5E+004 N	2.1E+002 N	5.4E+002 N	8.2E+005 N	3.1E+004 N	3.3E+000	6.5E+001 N
ETHYL ACETATE	141786	9.00E-001 I					5.5E+003 N	3.3E+003 N	1.2E+003 N	1.8E+006 N	7.0E+004 N	1.7E+000	3.5E+001 N
ETHYL BENZENE	100414	1.00E-001 I		2.90E-001 I		y	1.3E+003 N	1.1E+003 N	1.4E+002 N	2.0E+005 N	7.8E+003 N	7.5E-001	1.5E+001 N
ETHYLENE DIAMINE	107153	2.00E-002 H					7.3E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N		
ETHYLENE GLYCOL	107211	2.00E+000 I					7.3E+004 N	7.3E+003 N	2.7E+003 N	4.1E+006 N	1.6E+005 N	1.5E+001	
ETHYLENE GLYCOL, MONOBUTYL ETHER	111762	5.00E-001 I		3.70E+000 I			1.8E+004 N	1.4E+004 N	6.8E+002 N	1.0E+006 N	3.9E+004 N	1.5E+001	3.0E+002 N
ETHYLENE OXIDE	75218		1.00E+000 H		3.50E-001 H	y	2.3E-002 C	1.8E-002 C	3.2E-003 C	5.7E+000 C	6.4E-001 C	4.8E-006	9.5E-005 C
ETHYLENE THIOUREA	96457	8.00E-005 I	1.1E-001 H				6.1E-001 C	5.7E-002 C	2.9E-002 C	5.2E+001 C	5.8E+000 C		
ETHYL ETHER	60297	2.00E-001 I					1.2E+003 N	7.3E+002 N	2.7E+002 N	4.1E+005 N	1.6E+004 N	4.2E-001	8.5E+000 N
ETHYL METHACRYLATE	97632	9.00E-002 H					5.5E+002 N	3.3E+002 N	1.2E+002 N	1.8E+005 N	7.0E+003 N	1.0E+000	2.1E+001 N
FENAMIPHOS	22224926	2.50E-004 I					9.1E+000 N	9.1E-001 N	3.4E-001 N	5.1E+002 N	2.0E+001 N	7.8E-003	1.6E-001 N
FLUOMETURON	2164172	1.30E-002 I					4.7E+002 N	4.7E+001 N	1.8E+001 N	2.7E+004 N	1.0E+003 N		
FLUORINE	7782414	6.00E-002 I					2.2E+003 N	2.2E+002 N	8.1E+001 N	1.2E+005 N	4.7E+003 N		
FOMESAFEN	72178020		1.90E-001 I				3.5E-001 C	3.3E-002 C	1.7E-002 C	3.0E+001 C	3.4E+000 C		
FONOFOSS	944229	2.00E-003 I					7.3E+001 N	7.3E+000 N	2.7E+000 N	4.1E+003 N	1.6E+002 N	1.8E-001	3.5E+000 N
FORMALDEHYDE	50000	2.00E-001 I			4.50E-002 I		7.3E+003 N	7.3E+000 N	2.7E+000 N	4.1E+003 N	1.6E+002 N	1.5E+000	3.0E+001 N
FORMIC ACID	64186	2.00E+000 H					7.3E+003 N	1.4E-001 C	2.7E+002 N	4.1E+005 N	1.6E+004 N	1.5E+000	3.0E+001 N
FURAN	110009	1.00E-003 I					7.3E+004 N	7.3E+003 N	2.7E+003 N	4.1E+006 N	1.6E+005 N		
FURAZOLIDONE	67458		3.80E+000 H			y	6.1E+000 N	3.7E+000 N	1.4E+000 N	2.0E+003 N	7.8E+001 N	1.5E-003	3.0E-002 N
FURFURAL	98011	3.00E-003 I		1.00E-002 A			1.8E-002 C	1.6E-003 C	8.3E-004 C	1.5E+000 C	1.7E-001 C		
GLYCIDALDEHYDE	765344	4.00E-004 I		2.90E-004 H			1.1E+002 N	3.7E+001 N	4.1E+000 N	6.1E+003 N	2.3E+002 N	2.3E-002	4.6E-001 N
GLYPHOSATE	1071836	1.00E-001 I					1.5E+001 N	1.1E+000 N	5.4E-001 N	8.2E+002 N	3.1E+001 N		
HEPTACHLOR	78448	5.00E-004 I	4.50E-000 I		4.50E+000 I		3.7E+003 N	3.7E+002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N	2.6E+001	5.3E+002 N
HEPTACHLOR EPOXIDE	1024573	1.30E-005 I	9.10E+000 I		9.10E+000 I		1.5E-002 C	1.4E-003 C	7.0E-004 C	1.3E+000 C	1.4E-001 C	4.2E-002	8.4E-001 C
HEXBROMOBENZENE	87821	2.00E-003 I					7.4E-003 C	6.9E-004 C	3.5E-004 C	6.3E-001 C	7.0E-002 C	1.2E-003	2.5E-002 C
HEXACHLOROBENZENE	118741	8.00E-004 I	1.60E+000 I		1.60E+000 I		7.3E+001 N	7.3E+000 N	2.7E+000 N	4.1E+003 N	1.6E+002 N		
HEXACHLOROBUTADIENE	87683	2.00E-004 H	7.80E-002 I		7.80E-002 I		4.2E-002 C	3.9E-003 C	2.0E-003 C	3.6E+000 C	4.0E-001 C	2.6E-003	5.2E-002 C
ALPHA-HCH	319846		6.30E+000 I		6.30E+000 I		8.6E-001 C	8.0E-002 C	4.0E-002 C	7.3E+001 C	8.2E+000 C	9.2E-002	1.8E+000 C
BETA-HCH	319846		1.80E+000 I		1.80E+000 I		1.1E-002 C	9.9E-004 C	5.0E-004 C	9.1E-001 C	1.0E-001 C	4.5E-005	8.9E-004 C
GAMMA-HCH (LINDANE)	58897	3.00E-004 I	1.30E+000 H		1.80E+000 I		3.7E-002 C	3.5E-003 C	1.8E-003 C	3.2E+000 C	3.5E-001 C	1.6E-004	3.1E-003 C
TECHNICAL HCH	608731						5.2E-002 C	4.8E-003 C	2.4E-003 C	4.4E+000 C	4.9E-001 C	2.2E-004	4.3E-003 C
HEXACHLOROCYCLOPENTADIENE	77474	7.00E-003 I		1.80E+000 I		1.80E+000 I	3.7E-002 C	3.5E-003 C	1.8E-003 C	3.2E+000 C	3.5E-001 C		
HEXACHLORODIBENZODIOXIN MIX	19408743		6.20E+003 I		2.00E-005 H		2.6E+002 N	7.3E-002 N	9.5E+000 N	1.4E+004 N	5.5E+002 N	1.0E+002	2.0E+003 N
							4.55E+003 I	1.1E-005 C	1.4E-006 C	5.1E-007 C	9.2E-004 C	1.0E-004 C	

Sources: I = IRIS, H = HEAST, A = HEAST Alternate, W = Withdrawn from IRIS or HEAST E = EPA-NCEA provisional value, O = other										Basis: C = Carcinogenic effects, N = Noncarcinogenic effects, 1 = RBC at HI of 0.1 < RBC-C						Region III SSLs	
Chemical	CAS	RfDo mg/kg/d	CSFo 1/mg/kg/d	RfDi mg/kg/d	CSFi 1/mg/kg/d	VOC	Tap water ug/l	Ambient air ug/m3	Fish mg/kg	Soil Industrial mg/kg	Residential mg/kg	DAF 1 mg/kg	DAF 20 mg/kg	Soil, for groundwater migration			
HEXACHLOROETHANE	87721	1.00E-003 I	1.40E-002 I		1.40E-002 I		4.8E+000 C I	4.5E-001 C I	2.3E-001 C I	4.1E+002 C I	4.6E+001 C I	1.8E-002	3.6E-001 C				
HEXACHLOROPHENNE	70304	3.00E-004 I					1.1E+001 N	1.1E+000 N	4.1E-001 N	6.1E+002 N	2.3E+001 N	1.0E+002	2.0E+003 N				
1,6-HEXAMETHYLENE DIISOCYANATE	822060			2.90E-006 I			1.1E-002 N										
HEXANE	110543	6.00E-002 H		5.71E-002 I			3.5E+002 N	2.1E+002 N	8.1E+001 N	1.2E+005 N	4.7E+003 N		6.9E-001	1.4E+001 N			
2-HEXANONE	591786	4.00E-002 E		1.4E-003 E			1.5E+003 N	5.1E+000 N	5.4E+001 N	8.2E+004 N	3.1E+003 N						
HEXAZINONE	51235042	3.30E-002 I					1.2E+003 N	1.2E+002 N	4.5E+001 N	6.7E+004 N	2.6E+003 N						
HMX	2691410	5.00E-002 I					1.8E+003 N	1.8E+002 N	6.8E+001 N	1.0E+005 N	3.9E+003 N						
HYDRAZINE	302012		3.00E+000 I			1.70E+001 I	2.2E-002 C	3.7E-004 C	1.1E-003 C	1.9E+000 C	2.1E-001 C						
HYDROGEN CHLORIDE	7647010			5.70E-003 I			2.1E+001 N										
HYDROGEN SULFIDE	7783064	3.00E-003 I		2.85E-004 I			1.1E+002 N	1.0E+000 N	4.1E+000 N	6.1E+003 N	2.3E+002 N						
HYDROQUINONE	123319	4.00E-002 H					1.5E+003 N	1.5E+002 N	5.4E+001 N	8.2E+004 N	3.1E+003 N						
IRON	7439896	3.00E-001 E					1.1E+004 N	1.1E+003 N	4.1E+002 N	6.1E+005 N	2.3E+004 N						
ISOBUTANOL	78831	3.00E-001 I					1.8E+003 N	1.1E+003 N	4.1E+002 N	6.1E+005 N	2.3E+004 N	5.9E-001	1.2E+001 N				
ISOPHORONE	78591	2.00E-001 I	9.50E-004 I				7.0E+001 C	6.6E+000 C	3.3E+000 C	6.0E+003 C	6.7E+002 C	2.1E-002	4.1E-001 C				
ISOPROPALIN	33820530	1.50E-002 I					5.5E+002 N	5.5E+001 N	2.0E+001 N	3.1E+004 N	1.2E+003 N						
ISOPROPYL METHYL PHOSPHONIC ACID	1832548	1.00E-001 I					3.7E+003 N	3.7E+002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N						
TETRAETHYLLEAD	78002	1.00E-007 I					3.7E-003 N	3.7E-004 N	1.4E-004 N	2.0E-001 N	7.8E-003 N	4.6E-005	9.2E-004 N				
LITHIUM	7439932	2.00E-002 E					7.3E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N						
MALATHION	121755	2.00E-002 I					7.3E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N	4.0E-001	8.1E+000 N				
MALEIC ANHYDRIDE	108316	1.00E-001 I					3.7E+003 N	3.7E+002 N	1.4E+002 N	2.0E+005 N	7.8E+003 N						
MANGANESE-NONFOOD	7439965	2.00E-002 I		1.43E-005 I			7.3E+002 N	5.2E-002 N	2.7E+001 N	4.1E+004 N	1.6E+003 N	4.8E+001	9.5E+002 N				
MANGANESE-FOOD	7439965	1.40E-001 I		1.43E-005 I			5.1E+003 N	5.2E-002 N	1.9E+002 N	2.9E+005 N	1.1E+004 N	3.3E+002	6.7E+003 N				
MEPHOSFOLAN	950107	9.00E-005 H					3.3E+000 N	3.3E-001 N	1.2E-001 N	1.8E+002 N	7.0E+000 N						
MEPIQUAT CHLORIDE	24307264	3.00E-002 I					1.1E+003 N	1.1E+002 N	4.1E+001 N	6.1E+004 N	2.3E+003 N						
MERCURIC CHLORIDE	7487947	3.00E-004 I					1.1E+001 N	1.1E+000 N	4.1E-001 N	6.1E+002 N	2.3E+001 N						
MERCURY (INORGANIC)	7439976		8.60E-005 I					3.1E-001 N									
METHYLMERCURY	22967926	1.00E-004 I					3.7E+000 N	3.7E-001 N	1.4E-001 N	2.0E+002 N	7.8E+000 N						
METHACRYLONITRILE	126987	1.00E-004 I		2.00E-004 A			1.0E+000 N	7.3E-001 N	1.4E-001 N	2.0E+002 N	7.8E+000 N	2.1E-004	4.2E-003 N				
METHANOL	67561	5.00E-001 I					1.8E+004 N	1.8E+003 N	6.8E+002 N	1.0E+006 N	3.9E+004 N	3.8E+000	7.5E+001 N				
METHIDATHION	950378	1.00E-003 I					3.7E+001 N	3.7E+000 N	1.4E+000 N	2.0E+003 N	7.8E+001 N						
METHOXYCHLOR	72435	5.00E-003 I					1.8E+002 N	1.8E+001 N	6.8E+000 N	1.0E+004 N	3.9E+002 N	1.5E+001	3.1E+002 N				
METHYL ACETATE	79209	1.00E+000 H					6.1E+003 N	3.7E+003 N	1.4E+003 N	2.0E+006 N	7.8E+004 N	1.2E+000	2.5E+001 N				
METHYL ACRYLATE	96333	3.00E-002 A					1.8E+002 N	1.1E+002 N	4.1E+001 N	6.1E+004 N	2.3E+003 N	5.0E-001	1.0E+001 N				
2-METHYLANILINE	95534		2.40E-001 H				2.8E-001 C	2.6E-002 C	1.3E-002 C	2.4E+001 C	2.7E+000 C	2.8E-004	5.7E-003 C				
4-(2-METHYL-4-CHLOROPHOENOXY) BUTYRIC ACID	94815	1.00E-002 I					3.7E+002 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N						
2-METHYL-4-CHLOROPHOENOXYACETIC ACID (MCPA)	94746	5.00E-004 I					1.8E+001 N	1.8E+000 N	6.8E-001 N	1.0E+003 N	3.9E+001 N						
2-(2-METHYL-4-CHLOROPHOENOXY)PROPIONIC ACID (MC	93652	1.00E-003 I					3.7E+001 N	3.7E+000 N	1.4E+000 N	2.0E+003 N	7.8E+001 N						
METHYLCYCLOXANE	108872		8.60E-001 H				6.3E+003 N	3.1E+003 N									
METHYLENE BROMIDE	74953	1.00E-002 A					6.1E+001 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N	1.5E-002	3.0E-001 N				
METHYLENE CHLORIDE	75092	6.00E-002 I	7.50E-003 I	8.60E-001 H	1.65E-003 I	y	4.1E+000 C	3.8E+000 C	4.2E-001 C	7.6E+002 C	8.5E+001 C	9.5E-004	1.9E-002 C				
4,4'-METHYLENE BIS(2-CHLOROANILINE)	101144	7.00E-004 H	1.30E-001 H		1.30E-001 H		5.2E-001 C	4.8E-002 C	2.4E-002 C	4.4E+001 C	4.9E+000 C						
4,4'-METHYLENE BIS(N,N'-DIMETHYL)ANILINE	101611		4.60E-002 I				1.5E+000 C	1.4E-001 C	6.9E-002 C	1.2E+002 C	1.4E+001 C						
4,4'-METHYLEDIPHENYL ISOCYANATE	101688			1.7E-004 I					6.2E-001 N								
METHYLETHYL KETONE (2-BUTANONE)	78933	6.00E-001 I		2.86E-001 I		y	1.9E+003 N	1.0E+003 N	8.1E+002 N	1.2E+006 N	4.7E+004 N	4.0E-001	7.9E+000 N				
METHYL HYDRAZINE	60344		1.10E+000 W				6.1E-002 C	5.7E-003 C	2.9E-003 C	5.2E+000 C	5.8E-001 C						

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E = EPA-NCEA registration number

Basic C = Carcinogenic effects N = Noncarcinogenic effects * = RBC at HI of 0.1 < RBC <							
Chemical	CAS	Risk-based concentrations					Region III SSLs Soil, for groundwater migration
		RfDo mg/kg/d	CSFo 1/mg/kg/d	RfDi mg/kg/d	CSFi 1/mg/kg/d	VOC	
METHYL ISOBUTYL KETONE (4-METHYL-2-PENTANONE)	108101	8.00E-002 H		2.00E-002 A			
METHYL METHACRYLATE	80626	1.40E+000 I		2.00E-001 I		y	
2-METHYL-5-NITROANILINE	99558		3.30E-002 H			y	
METHYL PARATHION	298000	2.50E-004 I					
2-METHYLPHENOL	95487	5.00E-002 I					
3-METHYLPHENOL	108394	5.00E-002 I					
4-METHYLPHENOL	106445	5.00E-003 H					
METHYLSTYRENE MIX	25013154	6.00E-003 A		1.00E-002 A		y	
ALPHA-METHYLSTYRENE	98839	7.00E-002 A				y	
METHYL TERT-BUTYL ETHER	1634044			8.57E-001 I		y	
METOLACHLOR (DUAL)	51218452	1.50E-001 I					
MIREX	2385855	2.00E-004 I					
MOLYBDENUM	7439987	5E-003 I					
MONOCHLORAMINE	10599903	1E-001 I		1.00E-001 H			
NALED	300765	2E-003 I					
NICKEL REFINERY DUST					8.4E-001 I		
NICKEL	7440020	2.00E-002 I					
NITRATE	14797558	1.60E+000 I					
NITRIC OXIDE	10102439	1.00E-001 W					
NITRITE	14797650	1.00E-001 I				y	
2-NITROANILINE	88744		5.70E-005 H				
NITROBENZENE	98953	5.00E-004 I		6.00E-004 A		y	
NITROFURANTOIN	67209	7.00E-002 H					
NITROFURAZONE	59870		1.50E+000 H				
NITROGEN DIOXIDE	10102440	1.00E+000 W					
NITROGLYCERIN	55630		1.4E-002 E			y	
4-NITROPHENOL	100027	8.00E-003 E					
2-NITROPROPANE	79469		5.70E-003 I	9.40E+000 H	y		
N-NITROSO-DI-N-BUTYLAMINE	924163		5.40E+000 I	5.60E+000 I	y		
N-NITROSODIETHANOLAMINE	1116547	2.80E+000 I					
N-NITROSODIETHYLAMINE	55185	1.50E+002 I		1.50E+002 I			
N-NITROSODIMETHYLAMINE	62759	5.10E+001 I		5.10E+001 I			
N-NITROSODIPHENYLAMINE	86306	4.90E-003 I					
N-NITROSODIPROPYLAMINE	621647	7.00E+000 I					
N-NITROSO-N-ETHYLUREA	759739	1.40E+002 H					
N-NITROSO-N-METHYLETHYLAMINE	10595956	2.20E+001 I					
N-NITROSOPIRROLIDINE	930552	2.10E+000 I		2.10E+000 I			
M-NITROTOLUENE	99081	2.00E-002 E					
O-NITROTOLUENE	88722	1.00E-002 H				y	
P-NITROTOLUENE	99990	1.00E-002 H				y	
NUSTAR	85509199	7.00E-004 I				y	
ORYZALIN	19044883	5.00E-002 I					
OXADIAZON	19666309	5.00E-003 I					
OXAMYL	23135220	2.50E-002 I					
OXYFLUORFEN	42874033	3.00E-003 I					

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Chemical	CAS	RfDo mg/kg/d	CSFo 1/mg/kg/d	RfDi mg/kg/d	CSFi 1/mg/kg/d	VOC	Tap water ug/l	Ambient air ug/m3	Fish mg/kg	Soil Industrial mg/kg	Residential mg/kg			
PARAQUAT DICHLORIDE	1910425	4.50E-003 I					1.6E+002 N	1.6E+001 N	6.1E+000 N	9.2E+003 N	3.5E+002 N			
PARATHION	56382	6.00E-003 H					2.2E+002 N	2.2E+001 N	8.1E+000 N	1.2E+004 N	4.7E+002 N	5.0E-001	1.0E+001 N	
PENTACHLOROBENZENE	608935	8.00E-004 I					2.9E+001 N	2.9E+000 N	1.1E+000 N	1.6E+003 N	6.3E+001 N	1.0E+000	2.0E+001 N	
PENTACHLORONITROBENZENE	82688	3.00E-003 I	2.80E-001 H				2.6E-001 C	2.4E-002 C	1.2E+001 C	2.2E+001 C	2.5E+000 C	4.1E-003	8.2E-002 C	
PENTACHLOROPHENOL	87865	3.00E-002 I	1.20E-001 I				5.6E-001 C	5.2E-002 C	2.6E-002 C	4.8E+001 C	5.3E+000 C			
PERMETHRIN	52645531	5.00E-002 I					1.8E+003 N	1.8E+002 N	6.8E+001 N	1.0E+005 N	3.9E+003 N	1.2E+002	2.4E+003 N	
PHENOL	108952	6.00E-001 I					2.2E+004 N	2.2E+003 N	8.1E+002 N	1.2E+006 N	4.7E+004 N	6.7E+000	1.3E+002 N	
M-PHENYLENEDIAMINE	108452	6.00E-003 I					2.2E+002 N	2.2E+001 N	8.1E+000 N	1.2E+004 N	4.7E+002 N	4.9E-002	9.8E-001 N	
O-PHENYLENEDIAMINE	95545		4.70E-002 H				1.4E+000 C	1.3E-001 C	6.7E-002 C	1.2E+002 C	1.4E+001 C			
P-PHENYLENEDIAMINE	106503	1.90E-001 H					6.9E+003 N	6.9E+002 N	2.6E+002 N	3.9E+005 N	1.5E+004 N			
2-PHENYLPHENOL	90437		1.90E-003 H				3.5E+001 C	3.3E+000 C	1.7E+000 C	3.0E+003 C	3.4E+002 C			
PHOSPHINE	7803512	3.00E-004 I					1.1E+001 N	3.1E-001 N	4.1E-001 N	6.1E+002 N	2.3E+001 N			
PHOSPHORIC ACID	7664382		8.60E-005 I				1.1E+001 N							
PHOSPHORUS (WHITE)	7723140	2.00E-005 I					7.3E-001 N	7.3E-002 N	2.7E-002 N	4.1E+001 N	1.6E+000 N			
P-PHTHALIC ACID	100210	1.00E+000 H					3.7E+004 N	3.7E+003 N	1.4E+003 N	2.0E+006 N	7.8E+004 N			
PHthalic Anhydride	85449	2.00E+000 I					7.3E+004 N	1.3E+002 N	2.7E+003 N	4.1E+006 N	1.6E+005 N	2.6E+001	5.2E+002 N	
POLYBROMINATED BIPHENYLS							7.5E-003 C	7.0E-004 C	3.5E-004 C	6.4E-001 C	7.2E-002 C			
POLYCHLORINATED BIPHENYLS	1336363						3.3E-002 C	3.1E-003 C	1.6E-003 C	2.9E+000 C	3.2E-001 C	2.1E-002	4.1E-001 C	
ACROCLOR-1016	12674112	7.00E-005 I	7.00E-002 I				9.6E-001 C	8.9E-002 C	4.5E-002 C	8.2E+001 C	5.5E+000 N	2.1E-001	4.2E+000 C	
ACROCLOR-1221	11104282		2.00E+000 I				3.3E-002 C	3.1E-003 C	1.6E-003 C	2.9E+000 C	3.2E-001 C			
ACROCLOR-1232	11141165		2.00E+000 I				3.3E-002 C	3.1E-003 C	1.6E-003 C	2.9E+000 C	3.2E-001 C			
ACROCLOR-1242	53469219		2.00E+000 I				3.3E-002 C	3.1E-003 C	1.6E-003 C	2.9E+000 C	3.2E-001 C			
ACROCLOR-1248	12672296		2.00E+000 I				3.3E-002 C	3.1E-003 C	1.6E-003 C	2.9E+000 C	3.2E-001 C			
ACROCLOR-1254	11097691	2.00E-005 I	2.00E+000 I				3.3E-002 C	3.1E-003 C	1.6E-003 C	2.9E+000 C	3.2E-001 C	5.4E-002	1.1E+000 C	
ACROCLOR-1260	11096825		2.00E+000 I				3.3E-002 C	3.1E-003 C	1.6E-003 C	2.9E+000 C	3.2E-001 C			
POLYCHLORINATED TERPHENYLS	61788338		4.50E+000 E				1.5E-002 C	1.4E-003 C	7.0E-004 C	1.3E+000 C	1.4E-001 C			
POLYNUCLEAR AROMATIC HYDROCARBONS														
ACENAPHTHENE	83329	6.00E-002 I					3.7E+002 N	2.2E+002 N	8.1E+001 N	1.2E+005 N	4.7E+003 N	5.2E+000	1.0E+002 N	
ANTHRACENE	120127	3.00E-001 I					1.8E+003 N	1.1E+003 N	4.1E+002 N	6.1E+005 N	2.3E+004 N	2.3E+001	4.7E+002 N	
BENZ[A]ANTHRACENE	56553		7.30E-001 E				9.2E-002 C	8.6E-003 C	4.3E-003 C	7.8E+000 C	8.7E-001 C	7.3E-002	1.5E+000 C	
BENZO[B]FLUORANTHENE	205992		7.30E-001 E				9.2E-002 C	8.6E-003 C	4.3E-003 C	7.8E+000 C	8.7E-001 C	2.3E-001	4.5E+000 C	
BENZO[K]FLUORANTHENE	207089		7.30E-002 E				9.2E-001 C	8.6E-002 C	4.3E-002 C	7.8E+001 C	8.7E+000 C	2.3E+000	4.5E+001 C	
BENZO[A]PYRENE	50328		7.30E+000 I				9.2E-003 C	2.0E-003 C	4.3E-004 C	7.8E-001 C	8.7E-002 C	1.9E-002	3.7E-001 C	
CARBAZOLE	86748		2.00E-002 H				3.3E+000 C	3.1E-001 C	1.6E-001 C	2.9E+002 C	3.2E+001 C	2.3E-002	4.7E-001 C	
CHRYSENE	218019		7.30E-003 E				9.2E+000 C	8.6E-001 C	4.3E-001 C	7.8E+002 C	8.7E+001 C	7.3E+000	1.5E+002 C	
DIBENZ[A,H]ANTHRACENE	53703		7.30E+000 E				9.2E-003 C	8.6E-004 C	4.3E-004 C	7.8E-001 C	8.7E-002 C	7.0E-002	1.4E+000 C	
DIBENZOFURAN	132649	4.00E-003 E					2.4E+001 N	1.5E+001 N	5.4E+000 N	8.2E+003 N	3.1E+002 N	3.8E-001	7.7E+000 N	
FLUORANTHENE	206440	4.00E-002 I					1.5E+003 N	1.5E+002 N	5.4E+001 N	8.2E+004 N	3.1E+003 N	3.1E+002	6.3E+003 N	
FLUORENE	86737	4.00E-002 I					2.4E+002 N	1.5E+002 N	5.4E+001 N	8.2E+004 N	3.1E+003 N	6.8E+000	1.4E+002 N	
INDENO[1,2,3-C,D]PYRENE	193395		7.30E-001 E				9.2E-002 C	8.6E-003 C	4.3E-003 C	7.8E+000 C	8.7E-001 C	6.4E-001	1.3E+001 C	
2-METHYLNAPHTHALENE	91576	2.00E-002 E					y	1.2E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N	1.1E+000	2.2E+001 N
NAPHTHALENE	91203	2.00E-002 I					y	6.5E+000 N	3.3E+000 N	2.7E+001 N	4.1E+004 N	1.6E+003 N	7.7E-003	1.5E-001 N
PYRENE	129000	3.00E-002 I					y	1.8E+002 N	1.1E+002 N	4.1E+001 N	6.1E+004 N	2.3E+003 N	3.4E+001	6.8E+002 N
PROMETON	1610180	1.50E-002 I						5.5E+002 N	5.5E+001 N	2.0E+001 N	3.1E+004 N	1.2E+003 N		
PROMETRYN	7287196	4.00E-003 I						1.5E+002 N	1.5E+001 N	5.4E+000 N	8.2E+003 N	3.1E+002 N		

Sources: I = IRIS; H = HEAST; A = HEAST Alternate; W = Withdrawn from IRIS or HEAST  
 E = EPA-NCEA provisional value; O = other

Chemical	CAS	RfDo mg/kg/d	CSFo 1/mg/kg/d	RfDi mg/kg/d	CSFi 1/mg/kg/d	VOC	Risk-based concentrations					Region III SSLS		
							Tap water ug/l	Ambient air ug/m3	Fish mg/kg	Soil Industrial mg/kg	Residential mg/kg	Soil, for groundwater migration DAF 1 mg/kg	DAF 20 mg/kg	
PROPACHLOR	1918167	1.30E-002 I					4.7E+002 N	4.7E+001 N	1.8E+001 N	2.7E+004 N	1.0E+003 N			
PROPANIL	709988	5.00E-003 I					1.8E+002 N	1.8E+001 N	6.8E+000 N	1.0E+004 N	3.9E+002 N			
PROPARGITE	2312358	2.00E-002 I					7.3E+002 N	7.3E+001 N	2.7E+001 N	4.1E+004 N	1.6E+003 N			
N-PROPYLBENZENE	103651	4.00E-002 E					2.4E+002 N	1.5E+002 N	5.4E+001 N	8.2E+004 N	3.1E+003 N			
PROPYLENE GLYCOL	57556	2.00E+001 H					7.3E+005 N	7.3E+004 N	2.7E+004 N	4.1E+007 N	1.6E+006 N			
PROPYLENE GLYCOL, MONOETHYL ETHER	52125538	7.00E-001 H					2.6E+004 N	2.6E+003 N	9.5E+002 N	1.4E+006 N	5.5E+004 N			
PURSUIT	107982	7.00E-001 H				5.70E-001 I	2.6E+004 N	2.1E+003 N	9.5E+002 N	1.4E+006 N	5.5E+004 N			
PYRIDINE	81335775	2.50E-001 I					9.1E+003 N	9.1E+002 N	3.4E+002 N	5.1E+005 N	2.0E+004 N			
QUINOLINE	110861	1.00E-003 I					3.7E+001 N	3.7E+000 N	1.4E+000 N	2.0E+003 N	7.8E+001 N			
RDX	91225		1.20E+001 H				5.6E-003 C	5.2E-004 C	2.6E-004 C	4.8E-001 C	5.3E-002 C			
RESMETHRIN	121824	3.00E-003 I	1.10E-001 I				6.1E-001 C	5.7E-002 C	2.9E-002 C	5.2E-001 C	5.8E+000 C			
RONNEL	10453868	3.00E-002 I					1.1E+003 N	1.1E+002 N	4.1E+001 N	6.1E+004 N	2.3E+003 N			
ROtenone	299843	5.00E-002 H					1.8E+003 N	1.8E+002 N	6.8E+001 N	1.0E+005 N	3.9E+003 N			
SELENIUS ACID	83794	4.00E-003 I					1.5E+002 N	1.5E+001 N	5.4E+000 N	8.2E+003 N	3.1E+002 N			
SELENIUM	7783008	5.00E-003 I					1.8E+002 N	1.8E+001 N	6.8E+000 N	1.0E+004 N	3.9E+002 N			
SILVER	7782492	5.00E-003 I					1.8E+002 N	1.8E+001 N	6.8E+000 N	1.0E+004 N	3.9E+002 N			
SIMAZINE	7440224	5.00E-003 I					1.8E+002 N	1.8E+001 N	6.8E+000 N	1.0E+004 N	3.9E+002 N	9.5E-001	1.9E+001 N	
SODIUM AZIDE	122349	5.00E-003 I	1.20E-001 H				5.6E-001 C	5.2E-002 C	2.6E-002 C	4.8E-001 C	5.3E+000 C	1.6E+000	3.1E+001 N	
SODIUM DIETHYLDITHIOCARBAMATE	26628228	4.00E-003 I					1.5E+002 N	1.5E+001 N	5.4E+000 N	8.2E+003 N	3.1E+002 N	1.7E-004	3.3E-003 C	
STRONTIUM, STABLE	148185	3.00E-002 I	2.70E-001 H				2.5E-001 C	2.3E-002 C	1.2E-002 C	2.1E+001 C	2.4E+000 C	7.7E+002	1.5E+004 N	
STRYCHNINE	7440246	6.00E-001 I					2.2E+004 N	2.2E+003 N	8.1E+002 N	1.2E+006 N	4.7E+004 N	8.3E-003	1.7E-001 N	
STYRENE	57249	3.00E-004 I					1.1E+001 N	1.1E+000 N	4.1E-001 N	6.1E+002 N	2.3E+001 N			
2,3,7,8-TETRACHLORODIBENZODIOXIN	100425	2.00E-001 I				2.86E-001 I	1.6E+003 N	1.0E+003 N	2.7E+002 N	4.1E+005 N	1.6E+004 N			
1,2,4,5-TETRACHLOROBENZENE	1746016		1.50E+005 H				4.5E-007 C	4.2E-008 C	2.1E-008 C	3.8E-005 C	4.3E-006 C	4.3E-007	8.6E-006 C	
1,1,1,2-TETRACHLOROETHANE	95943	3.00E-004 I					1.1E+001 N	1.1E+000 N	4.1E-001 N	6.1E+002 N	2.3E+001 N	3.3E-002	6.6E-001 N	
1,1,2-TETRACHLOROETHANE	630206	3.00E-002 I	2.60E-002 I				2.60E-002 I	4.1E-001 C	2.4E-001 C	1.2E-001 C	2.2E+002 C	2.5E+001 C	2.0E-004	4.0E-003 C
TETRACHLOROETHENE	79345	6.00E-002 E	2.00E-001 I				2.00E-001 I	5.3E-002 C	3.1E-002 C	1.6E-002 C	2.9E+001 C	3.2E+000 C	3.4E-005	6.8E-004 C
2,3,4,6-TETRACHLOROPHENOL	127184	1.00E-002 I	5.20E-002 E	1.4E-001 E	1.4E-001 E	2.00E-003 E	1.1E+000 C	3.1E+000 C	6.1E-002 C	1.1E+002 C	1.2E+001 C	2.4E-003	4.8E-002 C	
P,A,A-TETRACHLOROTOLUENE	58902	3.00E-002 I					1.1E+003 N	1.1E+002 N	4.1E+001 N	6.1E+004 N	2.3E+003 N			
1,1,1,2-TETRAFLUOROETHANE	5216251		2.00E+001 H				3.3E-003 C	3.1E-004 C	1.6E-004 C	2.9E-001 C	3.2E-002 C			
TETRAHYDROFURAN	811972						1.7E+005 N	8.4E+004 N						
TETRYL	109999	2.00E-001 E	7.6E-003 E	8.6E-002 E	6.8E-003 E		8.8E+000 C	9.2E-001 C	4.2E-001 C	7.5E+002 C	8.4E+001 C			
THALLIC OXIDE	479458	1.00E-002 H					3.7E+002 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N			
THALLIUM	1314325	7.00E-005 W					2.6E+000 N	2.6E-001 N	9.5E-002 N	1.4E+002 N	5.5E+000 N			
THALLIUM ACETATE	7440280	7.00E-005 O					2.6E+000 N	2.6E-001 N	9.5E-002 N	1.4E+002 N	5.5E+000 N	1.8E-001	3.6E+000 N	
THALLIUM CARBONATE	563688	9.00E-005 I					3.3E+000 N	3.3E-001 N	1.2E-001 N	1.8E+002 N	7.0E+000 N			
THALLIUM CHLORIDE	6533739	8.00E-005 I					2.9E+000 N	2.9E-001 N	1.1E-001 N	1.6E+002 N	6.3E+000 N			
THALLIUM NITRATE	7791120	8.00E-005 I					2.9E+000 N	2.9E-001 N	1.1E-001 N	1.6E+002 N	6.3E+000 N			
THALLIUM SULFATE (2.1)	10102451	9.00E-005 I					3.3E+000 N	3.3E-001 N	1.2E-001 N	1.8E+002 N	7.0E+000 N			
THIOBENCARB	7446186	8.00E-005 I					2.9E+000 N	2.9E-001 N	1.1E-001 N	1.6E+002 N	6.3E+000 N			
TIN	28249776	1.00E-002 I					3.7E+002 N	3.7E+001 N	1.4E+001 N	2.0E+004 N	7.8E+002 N			
	7440315	6.00E-001 H					2.2E+004 N	2.2E+003 N	8.1E+002 N	1.2E+006 N	4.7E+004 N			

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